

Package: rcosmo (via r-universe)

October 29, 2026

URL <https://github.com/frycast/rcosmo>

BugReports <https://github.com/frycast/rcosmo/issues>

Title Cosmic Microwave Background Data Analysis

Version 1.1.5

Description Tools for handling and analysing spherical data on a HEALPix grid, with applications to cosmic microwave background maps. Supports reading FITS files, converting between spherical and Cartesian coordinates and HEALPix ordering schemes, selecting sky regions, sampling pixels, and visualising maps. Implements empirical covariance, correlation and variogram estimation, covariance model fitting, angular power spectrum estimation, and distributional and entropy-based comparisons of spherical data. The package and its methods are described by Fryer, Li and Olenko (2020) <[doi:10.32614/RJ-2020-012](https://doi.org/10.32614/RJ-2020-012)>.

Depends R (>= 4.5.0)

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Encoding UTF-8

LazyData true

Imports FITSio (>= 2.1-0), Rcpp (>= 0.12.11), mmap (>= 0.6-17), tibble (>= 1.4.2), rgl (>= 0.99.16), cli (>= 1.0.0), entropy (>= 1.2.1), geoR (>= 1.7-5.2.1), nnls (>= 1.4)

Suggests testthat, gsl

LinkingTo Rcpp

Config/roxygen2/version 8.0.0

Config/pak/sysreqs cmake libfreetype6-dev libglu1-mesa-dev make texlive libpng-dev libuv1-dev libgl1-mesa-dev zlib1g-dev

Repository <https://frycast.r-universe.dev>

Date/Publication 2026-10-19 08:11:49 UTC

RemoteUrl <https://github.com/frycast/rcosmo>

RemoteRef HEAD

RemoteSha 633a566d751b787250894f278a5ebdaa25b735d1

Contents

ancestor	4
areCompatibleCMBDFs	5
as.CMBDataFrame	6
assumedConvex	7
assumedUniquePix	8
baseNeighbours	8
cbind.CMBDataFrame	9
chi2CMB	10
children	11
CMBDat	11
CMBDataFrame	12
CMBWindow	14
coords	15
coords.CMBDataFrame	15
coords.CMBWindow	16
coords.data.frame	17
coords.HPDataFrame	18
corrCMB	19
covCMB	21
covmodelCMB	22
covPwSp	24
displayPixelBoundaries	25
displayPixels	26
downloadCMBMap	28
downloadCMBPS	29
entropyCMB	30
exprob	31
extrCMB	32
fmf	32
fRen	33
geo2sph	34
geoAngle	35
geoArea	36
geoArea.CMBDataFrame	36
geoArea.CMBWindow	37
geoArea.HPDataFrame	38
geoDist	38
header	39
healpixCentered	40
HPDataFrame	40
ibp2p	42
is.CMBDat	43
is.CMBDataFrame	43
is.CMBWindow	44
is.HPDataFrame	45
linesCMB	45

maxDist	46
maxWindowDist	47
minDist	48
neighbours	49
nest2ring	50
nestSearch	51
nside	52
nside.CMBDataFrame	53
nside.HPDataFrame	53
numeric2col	54
ordering	55
ordering.CMBDataFrame	55
ordering.HPDataFrame	56
p2bp	57
p2ibp	57
parent	58
pix	58
pix.CMBDataFrame	59
pix.HPDataFrame	60
pix2coords	60
pixelArea	61
pixelWindow	62
plot.CMBCorrelation	62
plot.CMBCovariance	63
plot.CMBDataFrame	64
plot.CMBWindow	65
plot.HPDataFrame	67
plot.variogram	68
plotAngDis	69
plotcovmodelCMB	70
plotvariogram	71
practicalRangeCMB	72
print.CMBDataFrame	73
print.HPDataFrame	73
pwSpCorr	74
qqnormWin	75
qqplotWin	76
qstat	77
rbind.CMBDataFrame	78
rcosmo	79
resolution	80
ring2nest	80
sampleCMB	81
siblings	82
sphericalHarmonics	82
summary.CMBDataFrame	83
summary.CMBWindow	84
summary.HPDataFrame	85

triangulate	85
variofitCMB	86
variogramCMB	88
window	90
window.CMBDat	91
window.CMBDataFrame	92
window.data.frame	94
window.HPDataFrame	96
winType	97
Index	99

ancestor	<i>Return index of kth ancestor pixel</i>
----------	---

Description

Gives the pixel at resolution $j - k$ that contains p , where p is specified at resolution j (notice it does not depend on j).

Usage

ancestor(p, k)

Arguments

- p A pixel index specified in nested order.
- k A generation of an ancestor pixel.

Value

A numeric vector of nested pixel indices at resolution $j - k$, identifying the ancestors containing the input pixels.

Examples

ancestor(4,2)
ancestor(17,2)

areCompatibleCMBDFs	<i>Check compatibleness of CMBDataFrames</i>
---------------------	--

Description

Compare attributes to decide if two CMBDataFrames are compatible

Usage

```
areCompatibleCMBDFs(cmbdf1, cmbdf2, compare.pix = FALSE)
```

Arguments

cmbdf1	a CMBDataFrame
cmbdf2	a CMBDataFrame
compare.pix	A boolean. If TRUE then cmbdf1 and cmbdf2 must share the same pixel indices to be considered compatible

Details

If the CMBDataFrames do not have compatible attributes then a message is printed indicating the attributes that do not match. To suppress this use the [suppressMessages](#) function

Value

A single logical value, TRUE if both objects have the same nside and ordering, and also the same pixel indices when compare.pix = TRUE; otherwise FALSE. Incompatibilities are reported with a suppressible message.

Examples

```
a <- CMBDataFrame(nside = 2, ordering = "ring", coords = "cartesian")
b <- CMBDataFrame(nside = 1, ordering = "nested", coords = "spherical")
areCompatibleCMBDFs(a,b)

suppressMessages(areCompatibleCMBDFs(a,b))
```

as.CMBDataFrame	<i>Convert objects to CMBDataFrame</i>
-----------------	--

Description

Safely converts a `data.frame` to a `CMBDataFrame`. The rows of the `data.frame` are assumed to be in the HEALPix order given by `ordering`, and at the HEALPix resolution given by `nside`. Coordinates, if present, are assumed to correspond to HEALPix pixel centers. The coordinates must be named either `x,y,z` (cartesian) or `theta, phi` (spherical colatitude and longitude respectively). If `df` is a `HPDataFrame` then it is possible that `df` has attribute `healpixCentered = TRUE`, in which case `as.CMBDataFrame` will perform HEALPix centering of coordinates.

Usage

```
as.CMBDataFrame(df, ordering, nside, spix, drop.coords = FALSE)
```

Arguments

<code>df</code>	Any <code>data.frame</code> or <code>HPDataFrame</code> whose rows are in HEALPix order
<code>ordering</code>	character string that specifies the ordering scheme ("ring" or "nested")
<code>nside</code>	an integer 2^k that specifies the Nside (resolution) HEALPix parameter
<code>spix</code>	an integer vector that specifies the HEALPix pixel index corresponding to each row of <code>df</code> . If <code>spix</code> is left blank and <code>df</code> is a <code>data.frame</code> , then <code>df</code> is assumed to contain data for every pixel at resolution parameter <code>nside</code> (the full sky). In other words, in this case, the number of rows of <code>df</code> must be equal to $12 * \text{nside}^2$. However, if <code>spix</code> is left blank and <code>df</code> is a <code>CMBDataFrame</code> , then <code>spix</code> is set equal to <code>pix(df)</code>
<code>drop.coords</code>	A logical. If <code>df</code> is a <code>HPDataFrame</code> then it is possible that <code>df</code> has attribute <code>healpixCentered = TRUE</code> , in which case <code>as.CMBDataFrame</code> will perform HEALPix centering of coordinates. If <code>drop.coords = TRUE</code> then this will be done by dropping existing coordinates entirely (quicker) to rely only on HEALPix pixel indices. Otherwise if <code>drop.coords = FALSE</code> this will be done by replacing existing coordinates with locations of HEALPix pixel centers.

Value

A `CMBDataFrame`

Examples

```
## Example 1: Create df with no coords, then create CMBDataFrames cmbdf and
## df2 with spherical coords

df <- data.frame(I=rnorm(12))
df

cmbdf <- as.CMBDataFrame(df,ordering= "ring", nside=1)
```

```
summary(cmbdf)
pix(cmbdf)
coords(cmbdf)

df2 <- coords(cmbdf, new.coords = "spherical")
df2

## Example 2: Create CMBDataFrames for first 10 Healpix centers

df <- data.frame(I=rnorm(10))
df
cmbdf <- as.CMBDataFrame(df,ordering= "ring", nside=2, spix=1:10)
summary(cmbdf)
pix(cmbdf)
```

assumedConvex	Check if a CMBWindow is assumed convex.
---------------	---

Description

Initially any [CMBWindow](#) is not assumed convex. The assumedConvex attribute can be change for any [CMBWindow](#).

Usage

```
assumedConvex(win, assume.convex)
```

Arguments

win	a CMBWindow object
assume.convex	optionally change the assumedConvex attribute to TRUE or FALSE

Value

A logical value giving the convexity assumption when assume.convex is omitted; otherwise the [CMBWindow](#) with its convexity attribute updated.

Examples

```
win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
assumedConvex(win1)
win2 <- assumedConvex(win1, assume.convex = TRUE)
assumedConvex(win2)
assumedConvex(win1) <- TRUE
assumedConvex(win1)
```

assumedUniquePix	<i>Check if object was assumed to have unique HEALPix indices</i>
------------------	---

Description

The function checks object's attribute assumedUniquePix. The attribute is True if the object was assumed to have rows that correspond to unique HEALPix pixel indices.

Usage

```
assumedUniquePix(obj)
```

Arguments

obj Any object.

Value

A boolean. This is TRUE if obj is a [CMBDataFrame](#) or a [HPDataFrame](#) whose rows were assumed to correspond to unique HEALPix pixel indices.

Examples

```
hp1 <- HPDataFrame(I=rnorm(5), nside = 1, spix = c(1,1,2,2,3))
pix(hp1)
coords(hp1, new.coords = "cartesian")
assumedUniquePix(hp1)

sky <- CMBDataFrame(nside = 32, coords = "cartesian", ordering = "nested")
sky.s <- CMBDataFrame(sky, sample.size = 100)
hpdf <- HPDataFrame(sky.s, auto.spix = TRUE)
assumedUniquePix(hpdf)
```

baseNeighbours	<i>Return neighbours of base pixels</i>
----------------	---

Description

The base-resolution comprises twelve pixels. baseNeighbours returns a map from the base pixel index bp to the vector of base pixels that are neighbours of bp, in counterclockwise order of direction: S,SE,E,NE,N,NW,W,SW. The presence of -1 indicates that the corresponding direction is empty.

Usage

```
baseNeighbours(bp)
```

Arguments

bp The base pixel index

Value

A numeric vector of eight neighbouring base pixel indices in the order S, SE, E, NE, N, NW, W, SW. Missing neighbours are coded as -1. Returns NULL for an index outside 1 to 12.

Examples

```
## Return neighbours of base pixel 1
baseNeighbours(1)

## There is no base pixel 14, so baseNeighbours returns NULL
baseNeighbours(14)
```

cbind.CMBDataFrame [cbind](#) for CMBDataFrames

Description

Add a new column or columns (vector, matrix or data.frame) to a [CMBDataFrame](#). Note that method dispatch occurs on the first argument. So, the CMBDataFrame must be the first argument

Usage

```
## S3 method for class 'CMBDataFrame'
cbind(..., deparse.level = 1)
```

Arguments

... (generalized) vectors or matrices. Columns to bind.

deparse.level Integer controlling the construction of labels in the case of non-matrix-like arguments.

Details

See the documentation for [cbind](#)

Value

A CMBDataFrame containing the combined columns, with HEALPix metadata and pixel indices retained from the single CMBDataFrame input.

Examples

```
cmbdf <- CMBDataFrame(nside = 1, ordering = "nested", coords = "spherical")
cmbdf2 <- cbind(cmbdf, myData = rep(1, 12))
cmbdf2
```

chi2CMB

Chi-squared statistic for two CMBWindows

Description

This function returns the empirical chi-squared statistic for intensities observations from two [CMBWindows](#) of the specified [CMBDataFrame](#). The functions employs the function [chi2.empirical](#) and uses histogram counts of intensities for computations.

Usage

```
chi2CMB(cmbdf, win1, win2, intensities = "I")
```

Arguments

cmbdf	A CMBDataFrame .
win1	A CMBWindow
win2	A CMBWindow
intensities	A CMBDataFrame column with measured values.

Value

Estimated Chi-squared statistic for observations in two [CMBWindows](#). For small sample sizes and many zero counts Chi-squared statistic is inefficient.

References

[chi2.empirical](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
win1 <- CMBWindow(theta = c(0, pi/2, pi/2), phi = c(0, 0, pi/2))
win2 <- CMBWindow(theta = c(pi/2, pi, pi/2), phi = c(0, 0, pi/2))
chi2CMB(cmbdf, win1, win2)
```

children	<i>Return children of a pixel</i>
----------	-----------------------------------

Description

Gives four pixels at resolution $j + 1$ that are contained in p , where p is a pixel specified at resolution j (notice it does not depend on j).

Usage

```
children(p)
```

Arguments

`p` A pixel index specified in nested order.

Value

A numeric vector of four nested pixel indices at resolution $j + 1$, identifying the children of pixel p .

Examples

```
children(11)
```

CMBDat	<i>CMBDat class</i>
--------	---------------------

Description

The function CMBDat creates objects of class CMBDat. These are lists containing header information and other metadata as well as an element called data, whose columns may include, for example, the intensity (I), polarisation (Q, U), PMASK and TMASK. It also may contain an [mmap](#) object that points to the CMB map data table in the FITS file.

Usage

```
CMBDat(filename, mmap = FALSE, spix)
```

Arguments

`filename` The path to the fits file.
`mmap` A boolean indicating whether to use memory mapping.
`spix` The sample pixels (rows) to read from the FITS file binary data table (optional)

Value

A list containing header information and other metadata as well as an element called data where: If `mmap = FALSE` then a `data.frame` is included, named `data`, whose columns may include, for example, the intensity (I), polarisation (Q, U), PMASK and TMASK. If `mmap = TRUE` then a [mmap](#) object is returned that points to the CMB map data table in the FITS file.

Examples

```
filename <- system.file("extdata", "CMB_testmap.fits", package = "rcosmo")
cmbdat <- CMBDat(filename, spix = 1:10)
class(cmbdat)
cmbdat$nside
cmbdat$ordering
head(cmbdat$data)
```

CMBDataFrame

CMBDataFrame class

Description

The function `CMBDataFrame` creates objects of class `CMBDataFrame`. These are a special type of [data.frame](#) that carry metadata about, e.g., the HEALPix ordering scheme, coordinate system, and `nside` parameter.

Usage

```
CMBDataFrame(
  CMBData,
  coords,
  win,
  include.polar = FALSE,
  include.masks = FALSE,
  spix,
  sample.size,
  nside,
  ordering,
  I,
  ...
)
```

Arguments

CMBData	Can be a string location of FITS file, another <code>CMBDataFrame</code> , a <code>CMBDat</code> object, a HPDataFrame or unspecified.
coords	Can be "spherical," "cartesian", or unspecified (HEALPix only).

<code>win</code>	optional CMBWindow object that specifies a spherical polygon within which to subset the full sky CMB data.
<code>include.polar</code>	TRUE if polarisation data is required, otherwise FALSE.
<code>include.masks</code>	TRUE if TMASK and PMASK are required, otherwise FALSE.
<code>spix</code>	Optional vector of sample pixel indices, or a path to a file containing comma delimited sample pixel indices. The ordering scheme is given by <code>ordering</code> . If <code>ordering</code> is unspecified then CMBData must be either a CMBDataFrame or a FITS file and the ordering scheme is then assumed to match that of CMBData.
<code>sample.size</code>	If a positive integer is given, a simple random sample of size equal to <code>sample.size</code> will be taken from CMBData. If <code>spix</code> is specified then <code>sample.size</code> must be unspecified.
<code>nside</code>	Optionally specify the <code>nside</code> parameter manually $nside=2^k$ (usually 1024 or 2048).
<code>ordering</code>	Specifies the desired HEALPix ordering scheme ("ring" or "nested") for the output CMBDataFrame. If <code>ordering</code> is unspecified then the ordering scheme will be taken from the CMBData object. If ordering information cannot be found the default will be "nested". This parameter also specifies the ordering scheme of <code>spix</code> .
<code>I</code>	A vector of intensities to be included if CMBData is unspecified. Note that $\text{length}(I)$ must equal $12 * nside^2$ if either <code>spix</code> or <code>sample.size</code> are unspecified.
<code>...</code>	Optional names data columns of length <code>nrow(CMBData)</code> to add to the CMBDataFrame.

Value

A CMBDataFrame whose `row.names` attribute contains HEALPix indices.

Examples

```
## Construct a small synthetic sky map.
df <- CMBDataFrame(nside = 16, I = rnorm(12 * 16 ^ 2),
                   ordering = "nested")

## Take a random sample.
df.sample <- CMBDataFrame(df, sample.size = 80)
plot(df.sample)

## Select pixels by index.
df.subset <- CMBDataFrame(df, spix = c(2,4,6))

## Inspect the map and the selected pixel indices.
summary(df)

pix(df.subset)
```

CMBWindow

*CMBWindow class.***Description**

The function CMBWindow creates objects of class CMBWindow. It is either a polygon or a disc type.

Usage

```
CMBWindow(..., r, set.minus = FALSE, assume.convex = FALSE)
```

Arguments

<code>...</code>	these arguments are compulsory and must be labelled either x, y, z (cartesian) or theta, phi (spherical, colatitude and longitude respectively). Alternatively, a single data.frame may be passed in with columns labelled x, y, z or theta, phi.
<code>r</code>	if a disc type window is required then this specifies the radius of the disc
<code>set.minus</code>	when TRUE the window will be the unit sphere minus the window specified
<code>assume.convex</code>	when TRUE the window is assumed to be convex resulting in a faster computation time when the window is used with functions such as subWindow . This argument is irrelevant when the window is not a polygon

Details

If `r` is unspecified then the rows of `...` correspond to counter-clockwise ordered vertices defining a spherical polygon lying entirely within one open hemisphere on the unit sphere. Counter-clockwise is understood from the perspective outside the sphere, facing the hemisphere that contains the polygon, looking toward the origin. Note that there must be at least 3 rows (vertices) to define a polygon (we exclude bygones). On the other hand, if `r` is specified then `...` must specify just one row, and this row is taken to be the center of a disc of radius `r`

Value

A data frame of class CMBWindow containing vertex coordinates for a polygon, or centre coordinates and radius `r` for a disc. Attributes record the coordinate system, window type and convexity assumption.

Examples

```
win <- CMBWindow(theta = c(pi/2,pi/2,pi/3, pi/3), phi = c(0,pi/3,pi/3,0))
plot(win)

## Create a disc type window
win1<- CMBWindow(x=0,y=3/5,z=4/5,r=0.8, set.minus =TRUE)
plot(win1)
```

```
## Apply a disc type window to CMBDataFrame
cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian", ordering = "nested")
window(cmbdf) <- CMBWindow(x=0,y=3/5,z=4/5,r=0.8, set.minus =TRUE)
plot(cmbdf)
```

coords	<i>Coordinate conversion generic</i>
--------	--------------------------------------

Description

Detailed descriptions and examples can be found in documentation for specific coords functions [coords.CMBDataFrame](#), [coords.CMBWindow](#), [coords.HPDataFrame](#), [coords.data.frame](#)

Usage

```
coords(x, ...)
```

Arguments

x	An object.
...	Unused arguments.

Value

A character string naming the coordinate system when it is queried, or an object of the input class with converted coordinates when a new system is supplied. See the individual methods for supported classes.

See Also

[coords.CMBDataFrame](#) [coords.CMBWindow](#) [coords.HPDataFrame](#) [coords.data.frame](#)

coords.CMBDataFrame	<i>Coordinate system from a CMBDataFrame</i>
---------------------	--

Description

If `new.coords` is unspecified then this function returns the coordinate system used in the CMBDataFrame `cmbdf`. The coordinate system is either "cartesian" or "spherical". If a new coordinate system is specified, using e.g. `new.coords = "spherical"`, then this function instead returns a new CMBDataFrame whose coordinates are of the specified type. The original CMBDataFrame, `cmbdf`, is unaffected. If you would like to change `cmbdf` without creating a new variable, then use [coords<-CMBDataFrame](#) (see examples below).

Usage

```
## S3 method for class 'CMBDataFrame'
coords(x, new.coords, ...)
```

Arguments

x	A CMBDataFrame, cmbdf.
new.coords	Specifies the new coordinate system ("spherical" or "cartesian") if a change of coordinate system is desired.
...	Unused arguments.

Value

If new.coords is unspecified, then the name of the coordinate system of cmbdf is returned. Otherwise a new CMBDataFrame is returned equivalent to cmbdf but having the desired change of coordinates

Examples

```
## Create df with no coords, then create df2 with cartesian coords
df <- CMBDataFrame(nside = 16)
df
coords(df)
df2 <- coords(df, new.coords = "cartesian")
coords(df2)

## Change the coords of df directly (to spherical)
coords(df) <- "spherical"
coords(df)
```

coords.CMBWindow

Coordinate system from a [CMBWindow](#)

Description

This function returns the coordinate system used in a [CMBWindow](#). The coordinate system is either "cartesian" or "spherical"

Usage

```
## S3 method for class 'CMBWindow'
coords(x, new.coords, ...)
```

Arguments

x	a CMBWindow, win.
new.coords	specifies the new coordinate system ("spherical" or "cartesian") if a change of coordinate system is desired
...	Unused arguments.

Details

If a new coordinate system is specified, using e.g. new.coords = "spherical", the coordinate system of the CMBWindow will be converted

Value

If new.coords is unspecified, then the name of the coordinate system of win is returned. Otherwise a new CMBWindow is returned equivalent to win but having the desired change of coordinates

Examples

```
## Create win with sperical coords, then change it to win1 with cartesian coords
win <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
coords(win)
win1 <- coords(win, new.coords = "cartesian")
coords(win1)

## Change back to spherical coordinates

coords(win1) <- "spherical"
coords(win1)
```

coords.data.frame	<i>Create a new data.frame with a given coordinate system</i>
-------------------	---

Description

This does not affect the original object unless new coordinate system is directly assigned.

Usage

```
## S3 method for class 'data.frame'
coords(x, new.coords, ...)
```

Arguments

x	a data.frame with columns labelled x, y, z (for cartesian) or theta, phi (for spherical colatitude and longitude respectively)
new.coords	specifies the new coordinate system ("spherical" or "cartesian").
...	Unused arguments.

Value

A new data.frame whose coordinates are as specified by new.coords

Examples

```
## Create df with no coords, then create df2 with spherical coords
df <- data.frame(x = c(1,0,0), y = c(0,1,0), z = c(0,0,1))
df

df2 <- coords(df, new.coords = "spherical")
df2

## The function coords does not affect the original object.
## To change the coords assign a new value ("spherical or "cartesian")

coords(df, new.coords = "spherical")
df
coords(df) <- "spherical"
df
```

coords.HPDataFrame *Coordinate system from a [HPDataFrame](#)*

Description

Add or change coordinates in a [HPDataFrame](#). This does not affect the argument object hpdf. Instead it returns a new [HPDataFrame](#) with the desired coordinates. To change hpdf directly see [coords<- .HPDataFrame](#).

Usage

```
## S3 method for class 'HPDataFrame'
coords(x, new.coords, healpixCentered = FALSE, ...)
```

Arguments

x	a HPDataFrame, hpdf.
new.coords	specifies the new coordinate system ("spherical" or "cartesian")
healpixCentered	boolean. If TRUE then columns x,y,z or theta, phi will be ignored and removed if present. This forces the coordinates to be found from HEALPix pixel indices only. Then the HEALPixCentered attribute of hpdf will be set to TRUE.
...	Unused arguments.

Details

If columns exist labelled x,y,z (cartesian) or theta, phi (colatitude and longitude respectively), then these will be treated as the coordinates of hpdf and converted accordingly. If columns x,y,z or theta,phi are not present then the healpix pixel indices as given by pix(hpdf) are used for assigning coordinates.

Value

A `HPDataFrame` with columns x,y,z (cartesian) or theta, phi (colatitude and longitude respectively)

Examples

```
df <- HPDataFrame(I = rep(0,12), nside = 1)
coords(df, new.coords = "cartesian")
# Notice that df is unchanged
df

# Instead, change df directly
coords(df) <- "spherical"

## specify cartesian coordinates then convert to spherical
hp1 <- HPDataFrame(x = c(1,0,0), y = c(0,1,0), z = c(0,0,1),
                  nside = 1, auto.spix = TRUE)
hp1 <- coords(hp1, new.coords = "spherical")

## Instead, ignore/drop existing coordinates and use HEALPix only
hp2 <- HPDataFrame(x = c(1,0,0), y = c(0,1,0), z = c(0,0,1),
                  nside = 1, auto.spix = TRUE)
hp2 <- coords(hp1, new.coords = "spherical", healpixCentered = TRUE)
```

corrCMB

Sample correlation function

Description

This function provides an empirical correlation function for data in a `CMBDataFrame` or `data.frame`. It assumes that data are from a stationary spherical random field and the correlation depends only on a geodesic distance between locations. Output is a binned correlation.

Usage

```
corrCMB(
  cmbdf,
  num.bins = 10,
  sample.size,
  max.dist = pi,
  breaks,
```

```

    equiareal = TRUE,
    calc.max.dist = FALSE
  )

```

Arguments

<code>cmbdf</code>	is a CMBDataFrame or data.frame
<code>num.bins</code>	specifies the number of bins
<code>sample.size</code>	optionally specify the size of a simple random sample to take before calculating correlation. This may be useful if the full correlation computation is too slow.
<code>max.dist</code>	an optional number between 0 and pi specifying the maximum geodesic distance to use for calculating correlation. Only used if breaks are unspecified.
<code>breaks</code>	optionally specify the breaks manually using a vector giving the break points between cells. This vector has length <code>num.bins</code> since the last break point is taken as <code>max.dist</code> . If <code>equiareal = TRUE</code> then these breaks should be $\cos(r_i)$ where r_i are radii. If <code>equiareal = FALSE</code> then these breaks should be r_i .
<code>equiareal</code>	if TRUE then the bins have equal spherical area. If false then the bins have equal annular widths. Default is TRUE.
<code>calc.max.dist</code>	if TRUE then the <code>max.dist</code> will be calculated from the locations in <code>cmbdf</code> . Otherwise either <code>max.dist</code> must be specified or <code>max.dist</code> will default to pi.

Value

#' An object of the class `CMBCorrelation` that is a modification of [variog](#) from the package **geoR** with variogram values replaced by correlation.

The attribute "breaks" contains the break points used to create bins. The result has `num.bins + 1` values since the first value at distance 0 is not counted as a bin.

u a vector with distances.

v a vector with estimated correlation values at distances given in **u**.

n number of pairs in each bin

sd standard deviation of the values in each bin

bins.lim limits defining the interval spanned by each bin

ind.bin a logical vector indicating whether the number of pairs in each bin is greater or equal to the value in the argument `pairs.min`

var.mark variance of the data

beta.ols parameters of the mean part of the model fitted by ordinary least squares

output.type echoes the option argument

max.dist maximum distance between pairs allowed in the correlation calculations

n.data number of data

direction direction for which the correlation was computed

call the function call

References

geoR package, [variog](#), [variogramCMB](#), [covCMB](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
corcmb <- corrCMB(cmbdf, max.dist = pi/2, num.bins = 6)
corcmb
```

covCMB	<i>Sample covariance function</i>
--------	-----------------------------------

Description

This function provides an empirical covariance function for data in a [CMBDataFrame](#) or `data.frame`. It assumes that data are from a stationary spherical random field and the covariance depends only on a geodesic distance between locations. Output is a binned covariance.

Usage

```
covCMB(
  cmbdf,
  num.bins = 10,
  sample.size,
  max.dist = pi,
  breaks,
  equiareal = TRUE,
  calc.max.dist = FALSE
)
```

Arguments

<code>cmbdf</code>	is a CMBDataFrame or data.frame
<code>num.bins</code>	specifies the number of bins
<code>sample.size</code>	optionally specify the size of a simple random sample to take before calculating covariance. This may be useful if the full covariance computation is too slow.
<code>max.dist</code>	an optional number between 0 and pi specifying the maximum geodesic distance to use for calculating covariance. Only used if <code>breaks</code> are unspecified.
<code>breaks</code>	optionally specify the breaks manually using a vector giving the break points between cells. This vector has length <code>num.bins</code> since the last break point is taken as <code>max.dist</code> . If <code>equiareal = TRUE</code> then these breaks should be $\cos(r_i)$ where r_i are radii. If <code>equiareal = FALSE</code> then these breaks should be r_i .
<code>equiareal</code>	if TRUE then the bins have equal spherical area. If false then the bins have equal annular widths. Default is TRUE.
<code>calc.max.dist</code>	if TRUE then the <code>max.dist</code> will be calculated from the locations in <code>cmbdf</code> . Otherwise either <code>max.dist</code> must be specified or <code>max.dist</code> will default to pi.

Value

An object of the class `CMB Covariance` that is a modification of `variog` from the package **geoR** with variogram values replaced by covariances.

The attribute "breaks" contains the break points used to create bins. The result has `num.bins + 1` values since the first value, the sample variance, is not counted as a bin.

u a vector with distances.

v a vector with estimated covariance values at distances given in **u**.

n number of pairs in each bin

sd standard deviation of the values in each bin

bins.lim limits defining the interval spanned by each bin

ind.bin a logical vector indicating whether the number of pairs in each bin is greater or equal to the value in the argument `pairs.min`

var.mark variance of the data

beta.ols parameters of the mean part of the model fitted by ordinary least squares

output.type echoes the option argument

max.dist maximum distance between pairs allowed in the covariance calculations

n.data number of data

direction direction for which the covariance was computed

call the function call

References

geoR package, `variog`, `variogramCMB`, `corrCMB`

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
Cov <- covCMB(cmbdf, max.dist = pi/2, num.bins = 6)
Cov
```

Description

This function computes the covariances given the separation distance of locations. Options for different covariance functions on spheres are available. The function extends the function `cov.spatial` for additional covariance models on spheres.

Usage

```
covmodelCMB(
  obj,
  cov.model = "matern",
  cov.pars = stop("no cov.pars argument provided"),
  kappa = 0.5
)
```

Arguments

<code>obj</code>	Vector of distances between pairs of spatial locations.
<code>cov.model</code>	A type of the correlation function. Available choices are: "matern", "exponential", "spherical", "powered.exponential", "cauchy", "gencauchy", "pure.nugget", "askey", "c2wendland", "c4wendland", "sinpower", "multiquadric". Default is "matern"
<code>cov.pars</code>	A vector with two covariance parameters. The first parameter corresponds to the variance σ^2 . The second parameter corresponds to the range ϕ of the correlation function.
<code>kappa</code>	A smoothness parameter of the correlation function.

Details

The function returns the value of the covariance $C(h)$ at the distance h . The covariance function has the form

$$C(h) = \sigma^2 * \rho(h/\phi).$$

The parameters of the covariance are positive numbers σ^2 , ϕ and κ .

Expressions for the correlation functions which are not included in the package **geoR**:

askey

$$\rho(h/\phi) = (1 - h/\phi)^{\kappa}, \text{ if } h < \phi;$$

$$0, \text{ otherwise.}$$

c2wendland

$$\rho(h/\phi) = (1 + \kappa * h/\phi) * (1 - h/\phi)^{\kappa}, \text{ if } h < \phi;$$

$$0, \text{ otherwise.}$$

c4wendland

$$\rho(h/\phi) = (1 + \kappa * h/\phi + (\kappa^2 - 1) * (h/\phi)^2 / 3) * (1 - h/\phi)^{\kappa}, \text{ if } h < \phi;$$

$$0, \text{ otherwise.}$$

sinpower

$$\rho(h/\phi) = 1 - (\sin(h/(2\phi)))^{\kappa}$$

multiquadric

$$C(h) = (1 - \phi)^{(2*\kappa)} / (1 + \phi^2 - 2 * \phi * \cos(h))^{\kappa}, 0 < \phi < 1$$

Additional information can be found in the section Details in [cov.spatial](#).

Value

Values of a covariance function for the given distances.

References

geoR package, [cov.spatial](#)
T. Gneiting. Strictly and non-strictly positive definite functions on spheres. Bernoulli 19 (2013), no. 4, 1327-1349.

Examples

```
## Compute Askey variogram at x = pi/4

1 - covmodelCMB(pi/4, cov.pars = c(1, pi), kappa = 3, cov.model = "askey" )

## Plot of the Askey covariance function

v1.f <- function(x, ...) {covmodelCMB(x, ...)}
curve( v1.f(x, cov.pars = c(1, 0.03), kappa = 3, cov.model = "askey"),
  from = 0, to = 0.1, xlab = "distance", ylab = expression(cov(h)), lty = 2,
  main = "covariance")
```

covPwSp	<i>Covariance estimate via power spectra</i>
---------	--

Description

This function provides a covariance estimate using the values of the estimated power spectra.

Usage

```
covPwSp(PowerSpectra, Ns)
```

Arguments

- PowerSpectra A data frame which first column lists values of multipole moments and the second column gives the corresponding values of CMB power spectra.
- Ns A number of points in which the covariance estimate is computed on the interval [-1,1]

Value

A data frame which first column is 1-d grid starting at -1+1/Ns and finishing at 1 with the step 2/Ns. The second column is the values of estimated covariances on this grid.

References

Formula (2.1) in Baran A., Terdik G. Power spectrum estimation of spherical random fields based on covariances. *Annales Mathematicae et Informaticae* 44 (2015) pp. 15–22.

Power Spectra data are from Planck Legacy Archive <https://www.cosmos.esa.int/web/planck/pla>

Examples

```
if (requireNamespace("gsl", quietly = TRUE)) {
  spectrum <- data.frame(l = 1:5, power = exp(-(1:5)))
  Cov_est <- covPwSp(spectrum, Ns = 100)
  plot(acos(Cov_est$t), Cov_est$Estimated_Cov, type = "l",
       xlab = "Angular distance", ylab = "Estimated covariance")
}
```

displayPixelBoundaries

Plot HEALPix pixel boundaries

Description

Plot the HEALPix pixel boundaries at nside

Usage

```
displayPixelBoundaries(
  nside,
  eps = pi/90,
  col = "gray",
  lwd = 1,
  ordering,
  incl.labels = 1:(12 * nside^2),
  nums.col = col,
  nums.size = 1,
  font = 2,
  depth_test = "always",
  ...
)
```

Arguments

nside	the HEALPix nside parameter (integer number 2^k)
eps	controls the smoothness of the plot, smaller eps implies more samples
col	the colour of plotted boundary lines
lwd	the thickness of the plotted boundary lines

ordering	optionally specify an ordering scheme from which to plot HEALPix pixel numbers. Can be either "ring" or "nested"
incl.labels	If ordering is specified then this parameter sets the pixel indices that will be displayed (default is all indices at nside)
nums.col	specifies the colour of pixel numbers if ordering is specified
nums.size	specifies the size of pixel numbers if ordering is specified
font	A numeric font number from 1 to 5, used if ordering is specified
depth_test	The depth test to be applied. This controls how resistant the plotted object is to being obscured. See rgl.material
...	arguments passed to <code>rgl::plot3d</code>

Value

Produces a plot of the HEALPix pixel boundaries.

Examples

```
displayPixelBoundaries(1, eps = pi/90, col = "red")
displayPixelBoundaries(2, eps = pi/90, col = "green")
```

displayPixels	<i>Display the pixels and grandchildren</i>
---------------	---

Description

Display the pixels `spix` at resolution `j` by colouring in the grandchildren of `spix` at resolution `plot.j`

Usage

```
displayPixels(
  boundary.j,
  j,
  plot.j = 5,
  spix,
  boundary.col = "gray",
  boundary.lwd = 1,
  incl.labels = 1:(12 * 4^boundary.j),
  col = "blue",
  size = 3
)
```

Arguments

boundary.j	The resolution to display boundaries at. If this is missing then boundaries will not be plotted.
j	The resolution that spix are specified at.
plot.j	The resolution to plot grandchildren at
spix	Integer vector. The pixel indices to display. These must be in nested order.
boundary.col	The boundary colour.
boundary.lwd	The boundary line width.
incl.labels	Integer vector of pixel indices to label at resolution j.
col	The colour to make the grandchildren.
size	The size to make the grandchildren.

Value

No meaningful return value; called for its side effect of drawing pixels and optional boundaries in the current rgl scene.

Examples

```
## Example 1

## Plot base pixels 1,2,3 by colouring their grandchildren at resolution
## 5 (by default). No pixel boundaries.
displayPixels(j=0, spix=c(1,2,3))

## Plot base pixels 1,2,3 display and their boundaries (boundary.j=0)
displayPixels(0,0, spix=c(1,2,3))

## Plot base pixels 1,2,3 by colouring their grandchildren at resolution 2
displayPixels(0,0, plot.j = 2, spix=c(1,2,3))

## Example 2

demoNeighbours <- function(p,j) {
  neighbours(p, j)
  displayPixels(boundary.j = j, j = j, plot.j = 5,
    spix = neighbours(p, j),
    boundary.col = "gray",
    boundary.lwd = 1,
    incl.labels = neighbours(p, j),
    col = "blue",
    size = 3)
  rcosmo::displayPixelBoundaries(nside = 1, col = "blue", lwd = 3)
}

demoNeighbours(1,2)
```

downloadCMBMap

Download CMB Maps from Planck Public Data Release.

Description

The function downloadCMBMap downloads CMB maps from https://irsa.ipac.caltech.edu/data/Planck/release_2/all-sky-maps/matrix_cmb.html.

Usage

```
downloadCMBMap(foreground = "smica", nside = 1024, destfile, release = 2)
```

Arguments

foreground	A string naming the foreground separation method pipeline. Please choose one of "COMMANDER", "NILC", "SEVEM" or "SMICA" (not case sensitive).
nside	An integer. The nside parameter (resolution) required. The available options are 1024 or 2048.
destfile	A required character string with an explicit path and file name for the downloaded file. No default destination is used. Tilde-expansion is performed.
release	An integer. Indicates the Planck map release to download. Currently only 2 and 3 are supported.

Details

CMB maps have been produced by the COMMANDER, NILC, SEVEM, and SMICA pipelines, respectively.

For each pipeline, the intensity maps are provided at Nside = 2048, at 5 arcmin resolution, and the polarization maps are provided at Nside = 1024 at 10 arcmin resolution.

Value

The integer status code from `utils::download.file`, returned invisibly (0 indicates success). The FITS map is saved to `destfile` and can be read with `CMBDataFrame`.

References

Planck Public Data Release 2 Maps https://irsa.ipac.caltech.edu/data/Planck/release_2/all-sky-maps/matrix_cmb.html

Other fits maps can also be downloaded using the general command [download.file](#).

Examples

```
## Requires internet access and downloads a large FITS file.
if (interactive()) {
  dest <- tempfile(fileext = ".fits")
  downloadCMBMap(foreground = "smica", nside = 1024, destfile = dest)
  sky <- CMBDataFrame(dest, spix = 1:10)
  unlink(dest)
}
```

downloadCMBPS

Download CMB Power Spectra from Planck Legacy Archive.

Description

The function downloadCMBPS downloads CMB power spectra components from <https://www.cosmos.esa.int/web/planck/pla>.

Usage

```
downloadCMBPS(link = 1, destfile, save = TRUE)
```

Arguments

link	The link code (an integer from 1 to 6) for the URL to download the file. See code details in this help file.
destfile	A character string with an explicit path and file name. Required when save = TRUE; unused when save = FALSE. No default destination is used. Tilde-expansion is performed.
save	A boolean indicating whether to save or not (since the downloaded data is returned anyway).

Details

link = 1: Best-fit LCDM CMB power spectra from the baseline Planck TT, TE, EE+lowE+lensing ($2 \leq \ell \leq 2508$).

link = 2: Baseline high- ℓ Planck TT power spectra ($2 \leq \ell \leq 2508$).

link = 3: Baseline high- ℓ Planck EE power spectra ($2 \leq \ell \leq 1996$).

link = 4: Baseline high- ℓ Planck TE power spectra ($2 \leq \ell \leq 1996$).

link = 5: Low- ℓ Planck EB power spectra ($2 \leq \ell \leq 29$).

link = 6: Low- ℓ Planck BB power spectra ($2 \leq \ell \leq 29$).

Value

A data frame returned invisibly, containing multipoles in column L and the selected power spectrum and uncertainty columns described above. When save = TRUE, it is also written to destfile.

References

Planck Legacy Archive <https://www.cosmos.esa.int/web/planck/pla>

Examples

```
## Requires internet access.
if (interactive()) {
  dest <- tempfile(fileext = ".txt")
  spectrum <- downloadCMBPS(link = 6, destfile = dest)
  unlink(dest)
  spectrum <- downloadCMBPS(link = 1, save = FALSE)
  plot(spectrum$L, spectrum$TT, type = "l")
}
```

entropyCMB

CMB Entropy

Description

This function returns an estimated entropy for the specified [CMBDataFrame](#) column intensities and [CMBWindow](#) region. The functions employs the function [entropy](#) and uses histogram counts of intensities for computations. All arguments of the standard [entropy](#) can be used.

Usage

```
entropyCMB(cmbdf, win, intensities = "I", method)
```

Arguments

cmbdf	A CMBDataFrame .
win	A CMBWindow
intensities	A CMBDataFrame column with measured values.
method	A method to estimate entropy, see entropy

Value

Estimated Shannon entropy for observations in [CMBWindow](#)

References

[entropy](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
win1 <- CMBWindow(theta = c(0, pi/2, pi/2), phi = c(0, 0, pi/2))
entropyCMB(cmbdf, win1)
```

exprob	<i>Threshold exceedance probability</i>
--------	---

Description

This function returns an estimated exceedance probability for the specified [CMBDataFrame](#) column intensities, threshold level *alpha* and [CMBWindow](#) region.

Usage

```
exprob(cmbdf, win, alpha, intensities = "I")
```

Arguments

cmbdf	A CMBDataFrame .
win	A CMBWindow
alpha	A numeric threshold level.
intensities	The name of the column in cmbdf that contains the measured values.

Value

Estimated threshold exceedance probability

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
win1 <- CMBWindow(theta = c(0, pi/2, pi/2), phi = c(0, 0, pi/2))
exprob(cmbdf, win1, alpha = mean(cmbdf$I))
```

extrCMB	<i>Extreme values</i>
---------	-----------------------

Description

This function returns n largest extreme values for the specified `CMBDataFrame` column intensities and `CMBWindow` region.

Usage

```
extrCMB(cmbdf, win, n, intensities = "I")
```

Arguments

<code>cmbdf</code>	A <code>CMBDataFrame</code> .
<code>win</code>	A <code>CMBWindow</code>
<code>n</code>	An integer value.
<code>intensities</code>	A <code>CMBDataFrame</code> column with measured values.

Value

A `CMBDataFrame` with n largest extreme values

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
win1 <- CMBWindow(theta = c(0, pi/2, pi/2), phi = c(0, 0, pi/2))
extremes <- extrCMB(cmbdf, win1, 5)
extremes
```

fmf	<i>First Minkowski functional</i>
-----	-----------------------------------

Description

This function returns an area of the spherical region where measured values are above of the specified threshold level *alpha*.

Usage

```
fmf(cmbdf, alpha, intensities = "I")
```

Arguments

cmbdf A [CMBDataFrame](#).
 alpha A numeric threshold level.
 intensities A [CMBDataFrame](#) column with measured values.

Value

The area of the exceedance region

References

Leonenko N., Olenko A. (2014) Sojourn measures of Student and Fisher-Snedecor random fields. Bernoulli, 20:1454-1483.

Examples

```
n <- 64
cmbdf <- CMBDataFrame(inside=n, I = rnorm(12*n^2),
                      coords = "cartesian",
                      ordering = "nested")

fmf(cmbdf, 0, 4)
fmf(cmbdf, 2, 4)

win <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
cmbdf.win <- window(cmbdf, new.window = win)
fmf(cmbdf.win, 0, 4)
```

fRen

Sample Renyi function

Description

This function computes values of the sample Renyi function. Returns the estimated values of $T(q)$ for q taking values on a grid. For large data sets could be rather time consuming.

Usage

```
fRen(
  cmbdf,
  q.min = 1.01,
  q.max = 10,
  N = 20,
  k.box = log2(inside(cmbdf)) - 3,
  intensities = "I"
)
```

Arguments

cmbdf	A CMBDataFrame .
q.min	Left endpoint of the interval to compute the Renyi function. The default value is 1.01,
q.max	Right endpoint of the interval to compute the Renyi function. The default value is 10
N	Number of points to compute the Renyi function. The default value is 20.
k.box	A dyadic decomposition level in computing the Renyi function, see the references in Details. The default value is $\log_2(\text{nside}(\text{cmbdf})) - 3$
intensities	A CMBDataFrame column with measured values

Value

Data frame which first column is the sampling grid $\text{seq}(q.\text{min}, q.\text{max}, \text{length.out} = N)$ of q values. Another column consists of values of the sample Renyi function $T(q)$ computed on the grid using the $k.\text{box}$ th level dyadic decomposition of the unit ball.

References

- (1) Leonenko, N., and Shieh, N. 2013. Rényi function for multifractal random fields. *Fractals* 21, Article No. 1350009.
- (2) <http://mathworld.wolfram.com/RenyiEntropy.html>

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
Tq <- fRen(cmbdf, N = 5, k.box = 1)
plot(Tq$q, Tq$Tq, xlab = "q", ylab = "Renyi function")
```

 geo2sph

Convert geographic to spherical coordinates

Description

Convert latitude (lat) and longitude (lon) to spherical coordinates (theta, phi) with theta in $[0, \pi]$ and phi in $[0, 2 * \pi]$. All values lat, lon, theta, phi are assumed to be in radians.

Usage

```
geo2sph(...)
```

Arguments

... A data.frame with columns lat and lon, or named vectors of lat and lon.

Value

A data.frame with columns theta and phi.

Examples

```
geo <- data.frame( lat = c(0, pi/3, pi/2), lon = c(0, pi/3, pi))
geo
sph <- geo2sph(geo)
sph
```

geoAngle	<i>Angle between two spherical directions</i>
----------	---

Description

Get an angle between two directions defined by arcs on the unit sphere

Usage

```
geoAngle(p1)
```

Arguments

p1 A [data.frame](#) with rows specifying numeric points located on the unit sphere. It should have columns labelled x,y,z for Cartesian or theta, phi for spherical colatitude and longitude respectively.

Value

Let $p1[1,]$, $p1[2,]$, and $p1[3,]$ denote the rows of p1. Then the returned object is an angle in radians between two arcs determined by # the pairs # of spherical points ($p1[1,]$, $p1[2,]$) and ($p1[2,]$, $p1[3,]$) respectively.

Examples

```
p1 <- data.frame(diag(3))
p1
colnames(p1) <- c("x", "y", "z")
geoAngle(p1)

geo <- data.frame( lat = c(30, 0, 20), lon = c(30, 60, 10))*(pi/180)
geo
p2 <- geo2sph(geo)
p2
geoAngle(p2)
```

geoArea	<i>geoArea generic</i>
---------	------------------------

Description

Detailed descriptions and examples can be found in documentation for specific geoArea functions [geoArea.CMBDataFrame](#), [geoArea.HPDataFrame](#), [geoArea.CMBWindow](#)

Usage

```
geoArea(x)
```

Arguments

x	An object.
---	------------

Value

A numeric scalar giving the area covered by the object on the unit sphere, in steradians.

See Also

[geoArea.CMBDataFrame](#) [geoArea.HPDataFrame](#) [geoArea.CMBWindow](#)

geoArea.CMBDataFrame	<i>Geodesic area covered by a CMBDataFrame</i>
----------------------	--

Description

Gives the surface on the unit sphere that is encompassed by all pixels in cmbdf

Usage

```
## S3 method for class 'CMBDataFrame'
geoArea(x)
```

Arguments

x	a CMBDataFrame.
---	-----------------

Value

The sum of the areas of all pixels (rows) in x.

Examples

```
## At low resolution, a few data points can
## occupy a large pixel area, e.g.:

cmbdf <- CMBDataFrame(nside = 1, spix = c(1,2,3))
pix(cmbdf)

## The total number of Healpix points at nside=1 equals 12. As cmbdf has 3 Healpix
## it occupies  $\pi = 1/4 \times (\text{surface area of unit sphere})$ 

geoArea(cmbdf)
plot(cmbdf, size = 5, hp.boundaries = 1)
```

geoArea.CMBWindow	Geodesic area of a CMBWindow
-------------------	--

Description

Geodesic area of a [CMBWindow](#)

Usage

```
## S3 method for class 'CMBWindow'
geoArea(x)
```

Arguments

x A CMBWindow.

Value

The spherical area inside the CMBWindow x.

Examples

```
## A window that covers 1/8 of the unit sphere is constructed and its area is
##  $\pi/2 = 1/8 \times (\text{surface area of unit sphere})$ 

win <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
geoArea(win)
```

geoArea.HPDataFrame	<i>Geodesic area covered by a HPDataFrame</i>
---------------------	---

Description

Gives the surface on the unit sphere that is encompassed by all pixels in x.

Usage

```
## S3 method for class 'HPDataFrame'
geoArea(x)
```

Arguments

x A HPDataFrame.

Value

The sum of the areas of all pixels (rows) in x.

Examples

```
## Generate random I for HPDataFrame
hp1 <- HPDataFrame(I=rnorm(5), nside = 1, spix = c(1,1,2,2,3))
pix(hp1)

## The total number of Healpix points at nside=1 equals 12. As hp1 has five
## I values at 3 Healpix points, then the occupied area is
## pi = 1/4*(surface area of unit sphere)

geoArea(hp1)
plot(hp1, size = 5, hp.boundaries = 1)
```

geoDist	<i>Geodesic distance on the unit sphere</i>
---------	---

Description

Get geodesic distance between points on the unit sphere

Usage

```
geoDist(p1, p2, include.names = FALSE)
```

Arguments

- `p1` A [data.frame](#) with rows specifying numeric points located on the unit sphere. It should have columns labelled x,y,z for Cartesian or theta, phi for spherical colatitude and longitude respectively.
- `p2` Same as `p1`.
- `include.names` Boolean. If TRUE then the row and column names of the returned matrix will be taken from the points in `p1` and `p2` (see examples below).

Value

Let n denote the number of rows of `p1` and let m denote the number of rows of `p2`. Then the returned object is an n by m matrix whose entry in position ij is the geodesic distance from the i th row of `p1` to the j th row of `p2`.

Examples

```
p1 <- data.frame(diag(3))
colnames(p1) <- c("x", "y", "z")
p1
p2 <- data.frame(x=c(1,0), y=c(0,3/5), z=c(0,4/5))
p2
geoDist(p1, p2, include.names = FALSE)
```

header

Get the FITS headers from a [CMBDataFrame](#)

Description

Get the FITS headers from a [CMBDataFrame](#)

Usage

```
header(cmbdf)
```

Arguments

`cmbdf` a [CMBDataFrame](#).

Value

The FITS headers belonging to the FITS file from which `cmbdf` data was imported

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
header(cmbdf)
```

healpixCentered	<i>Check if object is assumed to have HEALPix centered coordinates</i>
-----------------	--

Description

The function checks object's attribute `healpixCentered`. The attribute is `TRUE` if the object was assumed to have rows that correspond to unique HEALPix pixel indices.

Usage

```
healpixCentered(obj)
```

Arguments

`obj` Any object.

Value

A boolean. This is `TRUE` if `obj` is a [CMBDataFrame](#) or a [HPDataFrame](#) whose coordinates were assumed to correspond to HEALPix pixel center locations.

Examples

```
hp1 <- HPDataFrame(I=rnorm(5), nside = 1, spix = c(1,1,2,2,3))
pix(hp1)
coords(hp1, new.coords = "cartesian")
healpixCentered(hp1)

sky <- CMBDataFrame(nside = 32, coords = "cartesian", ordering = "nested")
sky.s <- CMBDataFrame(sky, sample.size = 100)
hpdf <- HPDataFrame(sky.s, auto.spix = TRUE)
healpixCentered(hpdf)
```

HPDataFrame	<i>HPDataFrame class</i>
-------------	--------------------------

Description

HPDataFrames are a type of data.frame designed to carry data located on the unit sphere. Each row of a HPDataFrame is associated with a HEALPix pixel index. The HPDataFrame also holds an attribute called `nside` which stores the HEALPix Nside parameter (i.e., the resolution of the HEALPix grid that is being used). Unlike [CMBDataFrame](#), HPDataFrames may have repeated pixel indices. They are made this way so that multiple data points falling within a given pixel can be stored in different rows of any given HPDataFrame.

Usage

```
HPDataFrame(
  ...,
  nside,
  ordering = "nested",
  auto.spix = FALSE,
  spix,
  assumedUniquePix = FALSE,
  delete.duplicates = FALSE,
  save.dots = FALSE,
  save.duplicate.indices = FALSE
)
```

Arguments

<code>...</code>	Data. Can be named vectors or a data.frame. May include columns (x,y,z) or (theta, phi) representing Cartesian or spherical coordinates of points on the unit sphere.
<code>nside</code>	Integer number 2^k , the nside parameter, i.e, resolution. If nside is unspecified, then the an attempt is made to use columns x,y and z from the provided data, as Cartesian coordinates, to calculate an nside that is sufficient to ensure all points belong to unique pixels.
<code>ordering</code>	The HEALPix ordering scheme ("ring" or "nested").
<code>auto.spix</code>	Boolean. If TRUE then spix will be found from the coordinates provided in the data. That is, each row of data will be assigned the pixel index of its closest HEALPix pixel center. There must be columns x,y,z for cartesian or theta, phi for spherical colatitude and longitude respectively. If <code>auto.spix = FALSE</code> then nside must be specified.
<code>spix</code>	A vector of HEALPix pixel indices indicating the pixel locations of the data. Note that spix is ignored if <code>auto.spix = TRUE</code> .
<code>assumedUniquePix</code>	A boolean. Sets the assumedUniquePix attribute of the HPDataFrame. This attribute indicates whether or not the rows of a HPDataFrame can be assumed to belong to unique pixels.
<code>delete.duplicates</code>	Boolean. If TRUE then rows corresponding to duplicate pixel indices will be dropped from the returned HPDataFrame, and assumedUniquePix will be set to TRUE.
<code>save.dots</code>	A logical. If TRUE then the dot product of each observation with the nearest child HEALPix pixel center will be stored as a column called "distance" in the returned HPDataFrame, provided that <code>auto.spix = TRUE</code> . Note that a 'child' pixel is any one of the four pixels contained in the current pixel, in the nested scheme, at the next highest resolution. See children .
<code>save.duplicate.indices</code>	A logical. If TRUE and <code>delete.duplicates</code> is also TRUE, then the row indices of duplicated pixels will be retained as an attribute called "duplicates". Note

that row index refers to the row position of the duplicated pixel in the original HPDataFrame, and not the actual pixel index itself.

Details

HPDataFrame with `auto.spix = TRUE` can be used to transform any spherical data (not necessarily CMB) to the Healpix representation, see Example 3 below.

Value

A data frame of class HPDataFrame containing the supplied observations and any coordinate columns. HEALPix pixel indices are stored in the `pix` attribute; further attributes record `nside`, ordering and whether indices are assumed unique.

Examples

```
hp <- HPDataFrame(I = c(1, 2, 3), nside = 1, spix = c(1, 1, 2))
pix(hp)
assumedUniquePix(hp)
coords(hp, new.coords = "cartesian")

## Assign pixels automatically from three distinct locations.
hp.unique <- HPDataFrame(x = c(1, 0, 0), y = c(0, 1, 0), z = c(0, 0, 1),
                        auto.spix = TRUE)

pix(hp.unique)
assumedUniquePix(hp.unique)
```

ibp2p

Computes pixel's index using its subindex within base resolution

Description

Find the pixel index p of a given pixel with index `ibp` in base pixel `bp`.

Usage

```
ibp2p(ibp, bp, j)
```

Arguments

<code>ibp</code>	The pixel index within base pixel <code>bp</code> , at resolution <code>j</code> , in nested order.
<code>bp</code>	The base pixel index
<code>j</code>	The resolution parameter $nside = 2^j$

Value

A numeric vector of full-sky nested pixel indices at resolution `j`, combining the within-base indices and base pixel indices.

Examples

```
ibp2p(1, 1, 2)
ibp2p(1, 2, 2)
```

is.CMBDat	<i>Check if an object is of class CMBDat</i>
-----------	--

Description

Check if an object is of class CMBDat

Usage

```
is.CMBDat(cmbdf)
```

Arguments

cmbdf	Any R object
-------	--------------

Value

TRUE if cmbdf is a CMBDat object, otherwise FALSE

Examples

```
filename <- system.file("extdata", "CMB_testmap.fits", package = "rcosmo")
cmbdat <- CMBDat(filename, spix = 1:10)
is.CMBDat(cmbdat)
```

is.CMBDataFrame	<i>Check if an object is of class CMBDataFrame</i>
-----------------	--

Description

Check if an object is of class CMBDataFrame

Usage

```
is.CMBDataFrame(cmbdf)
```

Arguments

cmbdf	Any R object
-------	--------------

Value

TRUE if cmbdf is a CMBDataFrame, otherwise FALSE

Examples

```
df <- CMBDataFrame(nside = 16)
is.CMBDataFrame(df)
df2 <- coords(df, new.coords = "cartesian")
is.CMBDataFrame(df2)
```

is.CMBWindow

Check if an object is a CMBWindow

Description

Check if an object is a CMBWindow

Usage

```
is.CMBWindow(win)
```

Arguments

win any object

Value

TRUE or FALSE depending if win is a CMBWindow

Examples

```
win <- CMBWindow(x=0,y=3/5,z=4/5,r=0.8, set.minus = TRUE)
is.CMBWindow(win)
```

is.HPDataFrame	Check if an object is of class HPDataFrame
----------------	--

Description

Check if an object is of class [HPDataFrame](#)

Usage

```
is.HPDataFrame(hpdf)
```

Arguments

hpdf	Any R object
------	--------------

Value

TRUE if hpdf is a HPDataFrame, otherwise FALSE

Examples

```
df <- CMBDataFrame(nside = 16)
is.HPDataFrame(df)

df <- HPDataFrame(I = rep(0,12), nside = 1)
is.HPDataFrame(df)
```

linesCMB	Adds lines of fitted variograms to variogram plots
----------	--

Description

This function adds a line with the variogram model fitted by the function [variofitCMB](#) to a current variogram plot. The function modifies [lines.variomodel.variofit](#) from the package **geoR** for additional covariance models on spheres.

Usage

```
linesCMB(x, max.dist, scaled = FALSE, ...)
```

Arguments

x	An object of the class variofit containing information about the fitted model obtained as an output of the function variofitCMB .
max.dist	A maximum distance to draw the variogram line. The default is <code>x\$max.dist</code> .
scaled	logical. If TRUE the sill in the plot is 1.
...	other plotting parameters passed to curve

Details

The function adds a line with fitted variogram model to a plot. It is used to compare empirical variograms against fitted models returned by [variofitCMB](#).

Available models are: "matern", "exponential", "spherical", "powered.exponential", "cauchy", "gen-cauchy", "pure.nugget", "askey", "c2wendland", "c4wendland", "sinpower", "multiquadric".

Value

A line with a fitted variogram model is added to a plot.

References

[geoR](#) package, [lines.variomodel.variofit](#), [covmodelCMB](#), [variofitCMB](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
varcmb <- variogramCMB(cmbdf, max.dist = pi/2, num.bins = 6)
ols <- variofitCMB(varcmb, ini.cov.pars = c(1, 0.5),
                  cov.model = "matern", weights = "equal", messages = FALSE)
plot(varcmb)
linesCMB(ols, lty = 2)
```

maxDist

Get the maximum geodesic distance between points

Description

Get the maximum geodesic distance either between all points in a data.frame pairwise, or between all points in a data.frame and one target point.

Usage

```
maxDist(df, point)
```

Arguments

df	A data.frame with columns x,y,z for cartesian or theta, phi for spherical colatitude and longitude respectively. The rows must correspond to points on the unit sphere. If this is a HPDataFrame or CMBDataFrame and coordinate columns are missing, then coordinates will be assigned based on HEALPix pixel indices.
point	An optional target point on the unit sphere in cartesian coordinates, in which case all distances are calculated between point and the points in df.

Value

If point is specified: the longest geodesic distance from point to the points specified by the rows of df. If point is not specified: the longest geodesic distance pairwise between points in df.

Examples

```
## Using a CMBDataFrame with HEALPix coordinates only
cmbdf <- CMBDataFrame(nside = 1, spix = c(1,5,12), ordering = "ring")
plot(cmbdf, hp.boundaries = 1, col = "blue", size = 5)
p <- c(0,0,1)
maxDist(cmbdf, p) # no need to have coordinates

## Using a HPDataFrame with HEALPix coordinates only
hp <- HPDataFrame(nside = 1, I = rep(0,3), spix = c(1,5,12) )
maxDist(hp, p) # notice no need to have coordinates

## Using a data.frame with cartesian coordinates
coords(hp) <- "cartesian"
df <- data.frame(x = hp$x, y = hp$y, z = hp$z)
maxDist(df, p)

## Using a data.frame with spherical coordinates
coords(hp) <- "spherical"
df <- data.frame(theta = hp$theta, phi = hp$phi)
maxDist(df, p)

## max distance between points in cmbdf
maxDist(cmbdf)
```

maxWindowDist

Get the maximum distance between all points in a [CMBWindow](#)

Description

Get the maximum distance between all points in a [CMBWindow](#)

Usage

```
maxWindowDist(x)
```

Arguments

x A CMBWindow object.

Value

The maximum distance between window's points.

Examples

```
## win is a equilateral spherical triangle which sides are pi/2
win <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
maxWindowDist(win)
```

minDist

Get the minimum geodesic distance between points

Description

Get the minimum geodesic distance either between all points in a data.frame pairwise, or between all points in a data.frame and one target point.

Usage

```
minDist(df, point)
```

Arguments

df	A data.frame with columns x,y,z for cartesian or theta, phi for spherical colatitude and longitude respectively. The rows must correspond to points on the unit sphere. If this is a HPDataFrame or CMBDataFrame and coordinate columns are missing, then coordinates will be assigned based on HEALPix pixel indices.
point	An optional target point on the unit sphere in cartesian coordinates, in which case all distances are calculated between point and the points in df.

Value

If point is specified: the shortest distance from point to the points specified by the rows of df. If point is not specified: the shortest distance pairwise between points in df.

Examples

```
## Using a CMBDataFrame with HEALPix coordinates only
cmbdf <- CMBDataFrame(nside = 1, spix = c(1,5,12), ordering = "ring")
plot(cmbdf, hp.boundaries = 1, col = "blue", size = 5)
p <- c(0,0,1)
minDist(cmbdf, p) # no need to have coordinates

## Using a HPDataFrame with HEALPix coordinates only
hp <- HPDataFrame(nside = 1, I = rep(0,3), spix = c(1,5,12) )
minDist(hp, p) # notice no need to have coordinates

## Using a data.frame with cartesian coordinates
coords(hp) <- "cartesian"
df <- data.frame(x = hp$x, y = hp$y, z = hp$z)
minDist(df, p)
```

```
## Using a data.frame with spherical coordinates
coords(hp) <- "spherical"
df <- data.frame(theta = hp$theta, phi = hp$phi)
minDist(df, p)

## min distance between points in cmdf
minDist(cmbdf)
```

neighbours	<i>Return neighbouring pixels</i>
------------	-----------------------------------

Description

Return the neighbouring pixels to a given pixel p that is specified at resolution j , in the nested order.

Usage

```
neighbours(p, j)
```

Arguments

p	Pixel index p at resolution j .
j	The resolution parameter with $n_{\text{side}} = 2^j$.

Value

A numeric vector of neighbouring pixel indices in nested order at resolution j ; directions without a neighbour are omitted. For $j > 0$, the vector also includes p itself. For a vector of input pixels, a list of these vectors is returned.

Examples

```
## Return the neighbouring pixels for base pixel 1
neighbours(1, 0)

## Plot the neighbouring pixels for base pixel 1
demoNeighbours <- function(p,j) {
  neighbours(p, j)
  displayPixels(boundary.j = j, j = j, plot.j = j+3,
    spix = neighbours(p, j),
    boundary.col = "gray",
    boundary.lwd = 1,
    incl.labels = neighbours(p, j),
    col = "blue",
    size = 3)
  rcosmo::displayPixelBoundaries(nside = 1, col = "blue", lwd = 3)
```

```

}

demoNeighbours(1,2)
demoNeighbours(1,0)

```

nest2ring

Convert nest to ring ordering

Description

Convert from "nested" to "ring" ordering

nest2ring computes the HEALPix pixel index in the "ring" ordering scheme from the pixel index in the "nested" ordering scheme.

Usage

```
nest2ring(nside, pix)
```

Arguments

nside	is the HEALPix nside parameter.
pix	is the set or subset of pixel indices at nside. If pix is left blank then all pixels are converted.

Value

the output is the corresponding set of pixel in the ring ordering scheme.

Examples

```

# compute HEALPix indices in the ring ordering scheme
nside <- 8
pix <-c(1,2,23)
nest2ring(nside,pix)

```

nestSearch	<i>Finds the closest pixel center to a point</i>
------------	--

Description

Finds the closest HEALPix pixel center to a given target point, specified in Cartesian coordinates, using an efficient nested search algorithm. HEALPix indices are all assumed to be in the "nested" ordering scheme.

Usage

```
nestSearch(target, nside, index.only = FALSE, save.dots = FALSE)
```

Arguments

target	A data.frame, matrix or vector of Cartesian (x,y,z) coordinates for the target point. If a data.frame is used then spherical coordinates can be specified with row names theta and phi.
nside	An integer, the target resolution at which the resulting pixels are returned.
index.only	A boolean indicating whether to return only the pixel index (TRUE), or cartesian coordinates as well (FALSE).
save.dots	A logical. A If TRUE then the dot product of each observation with the nearest child HEALPix pixel center will be returned as an attribute called "dot". Note that a 'child' pixel is any one of the four pixels contained in the current pixel in the nested scheme, at the next highest resolution. See children .

Value

if index.only = TRUE then the output will be a HEALPix index. If index.only FALSE then the output is the list containing the HEALPix index and Cartesian coordinate vector of the HEALPix point closest to target at resolution nside.

Examples

```
## Find the closest HEALPix pixel center at resolution j=2 for
## the point (0.6,0.8,0)

point <- c(0.6,0.8,0)
j <- 2
cpoint <- nestSearch(point, nside = 2^j)

## Plot the closest pixel center in blue and the point (0.6,0.8,0) in red
displayPixels(j, j, plot.j=j, spix=c(cpoint$pix),
              size=5, incl.labels =FALSE)
rgl::plot3d(point[1], point[2], point[3],
             col="red", size = 5, add = TRUE)
```

```
## Repeat the above for 4 points in a data.frame
points <- data.frame(x = c(1,0,0,0.6),
                    y = c(0,1,0,0.8),
                    z = c(0,0,1,0))

points
j <- 2
cpoints <- nestSearch(points, nside = 2^j)

## Plot the closest pixel center in blue and the point (0.6,0.8,0) in red
displayPixels(j, j, plot.j=j, spix=c(cpoints$pix),
             size=5, incl.labels =FALSE)
rgl::plot3d(points[,1], points[,2], points[,3],
            col="red", size = 5, add = TRUE)
```

nside

nside generic

Description

Detailed descriptions and examples can be found in documentation for specific nside functions [nside.CMBDataFrame](#), [nside.HPDataFrame](#)

Usage

```
nside(x)
```

Arguments

x An object.

Value

A numeric scalar giving the HEALPix nside parameter. The corresponding full-sky grid contains $12 \times \text{nside}^2$ pixels.

See Also

[nside.CMBDataFrame](#) [nside.HPDataFrame](#)

<code>nside.CMBDataFrame</code>	<i>HEALPix Nside parameter from a CMBDataFrame</i>
---------------------------------	--

Description

This function returns the HEALPix Nside parameter of a CMBDataFrame

Usage

```
## S3 method for class 'CMBDataFrame'
nside(x)
```

Arguments

`x` A [CMBDataFrame](#).

Value

The HEALPix Nside parameter.

Examples

```
df <- CMBDataFrame(nside = 16)
nside(df)
```

<code>nside.HPDataFrame</code>	<i>HEALPix Nside parameter from a HPDataFrame</i>
--------------------------------	---

Description

This function returns the HEALPix Nside parameter of a [HPDataFrame](#)

Usage

```
## S3 method for class 'HPDataFrame'
nside(x)
```

Arguments

`x` A [HPDataFrame](#).

Value

The HEALPix Nside parameter.

Examples

```
df <- HPDataFrame(I = rep(0,12), nside = 1)
nside(df)
```

numeric2col

numeric2col

Description

Map numeric values to a colour map

Usage

```
numeric2col(
  num,
  colmap = grDevices::terrain.colors(100),
  breaks.length = length(colmap)
)
```

Arguments

`num` A numeric vector. The numbers which will be mapped to colours.

`colmap` A colour map. See [palette](#).

`breaks.length` A single integer. Controls the number of breaks in the discretisation of `num`.

Value

A character vector of colours, one for each element of `num`, selected from `colmap` according to the discretised numeric values.

Examples

```
ns <- 16
sky <- CMBDataFrame(I = rnorm(12*ns^2), nside = ns)
plot(sky, col = numeric2col(sky$I))
```

ordering	<i>ordering generic</i>
----------	-------------------------

Description

Detailed descriptions and examples can be found in documentation for specific ordering functions [ordering.CMBDataFrame](#), [ordering.HPDataFrame](#)

Usage

```
ordering(x, ...)
```

Arguments

x	An object.
...	Extra arguments.

Value

A character string, "nested" or "ring", when queried, or an object of the input class reordered to the requested scheme.

See Also

[ordering.CMBDataFrame](#) [ordering.HPDataFrame](#)

ordering.CMBDataFrame	<i>HEALPix ordering scheme from a CMBDataFrame</i>
-----------------------	--

Description

This function returns the HEALPix ordering scheme from a CMBDataFrame. The ordering scheme is either "ring" or "nested".

Usage

```
## S3 method for class 'CMBDataFrame'
ordering(x, new.ordering, ...)
```

Arguments

x	A CMBDataFrame .
new.ordering	Specifies the new ordering ("ring" or "nest") if a change of ordering scheme is desired.
...	Unused arguments.

Details

If a new ordering is specified, using e.g. `new.ordering = "ring"`, the ordering scheme of the CMB-DataFrame will be converted.

Value

The name of the HEALPix ordering scheme that is used in the CMBDataFrame `x`.

Examples

```
df <- CMBDataFrame(nside = 1, ordering = "nested")
ordering(df)
df1 <- ordering(df, new.ordering = "ring")
ordering(df1)
```

ordering.HPDataFrame	<i>HEALPix ordering scheme from a HPDataFrame</i>
----------------------	---

Description

This function returns the HEALPix ordering scheme from a HPDataFrame. The ordering scheme is either "ring" or "nested". If a new ordering is specified, using e.g. `new.ordering = "ring"`, the ordering scheme of the HPDataFrame will be converted.

Usage

```
## S3 method for class 'HPDataFrame'
ordering(x, new.ordering, ...)
```

Arguments

<code>x</code>	a HPDataFrame .
<code>new.ordering</code>	Specifies the new ordering ("ring" or "nest") if a change of ordering scheme is desired.
<code>...</code>	Unused arguments.

Value

The name of the HEALPix ordering scheme that is used in the HPDataFrame `x`, or a new HP-DataFrame with the desired `new.ordering`

Examples

```
df <- HPDataFrame(I = rep(0,12), nside = 1, ordering = "nested")
ordering(df)
df1 <- ordering(df, new.ordering = "ring")
ordering(df1)
```

p2bp	<i>Return base pixel to which pixel belongs</i>
------	---

Description

The base pixel to which pixel p belongs at resolution j

Usage

p2bp(p, j)

Arguments

p	The pixel index at resolution j , in nested order.
j	The resolution parameter $n_{\text{side}} = 2^j$

Value

A numeric vector of base pixel indices from 1 to 12, identifying the base pixel containing each input pixel.

Examples

```
p2bp(5, 0)
p2bp(5, 1)
```

p2ibp	<i>Return pixel index within its base pixel</i>
-------	---

Description

Convert a pixel index p to its index within the base pixel to which p belongs

Usage

p2ibp(p, j)

Arguments

p	The pixel index at resolution j , in nested order.
j	The resolution parameter $n_{\text{side}} = 2^j$

Value

A numeric vector giving the nested index of each input pixel within its base pixel at resolution j .

Examples

```
p2ibp(6, 0)
p2ibp(6, 1)
```

parent	<i>Return index of parent pixel</i>
--------	-------------------------------------

Description

Gives the pixel at resolution $j - 1$ that contains p , where p is specified at resolution j (notice it does not depend on j).

Usage

```
parent(p)
```

Arguments

p A pixel index specified in nested order.

Value

A numeric vector of nested pixel indices at resolution $j - 1$, identifying the parent containing each input pixel.

Examples

```
parent(4)
parent(5)
```

pix	<i>pix generic</i>
-----	--------------------

Description

Detailed descriptions and and examples can be found in documentation for specific pix functions [pix.CMBDataFrame](#), [pix.HPDataFrame](#)

Usage

```
pix(x, ...)
```

Arguments

x	An object.
...	Extra arguments.

Value

A numeric vector of HEALPix pixel indices when queried, or an object of the input class with updated pixel indices when new indices are supplied.

See Also

[pix.CMBDataFrame](#) [pix.HPDataFrame](#)

pix.CMBDataFrame	<i>HEALPix pixel indices from</i> CMBDataFrame
------------------	--

Description

If new.pix is unspecified then this function returns the vector of HEALPix pixel indices from a CMBDataFrame. If new.pix is specified then this function returns a new CMBDataFrame with the same number of rows as cmbdf, but with pix attribute new.pix. Thus, new.pix must have length equal to nrow(cmbdf).

Usage

```
## S3 method for class 'CMBDataFrame'
pix(x, new.pix, ...)
```

Arguments

x	A CMBDataFrame.
new.pix	Optional vector of pixel indices with length equal to nrow(x).
...	Unused arguments.

Value

The vector of HEALPix pixel indices or, if new.pix is specified, a new CMBDataFrame.

Examples

```
df <- CMBDataFrame(nside = 16, sample.size = 10, ordering = "nested")
pix(df)
```

pix.HPDataFrame	<i>HEALPix pixel indices from HPDataFrame</i>
-----------------	---

Description

If new.pix is unspecified then this function returns the vector of HEALPix pixel indices from a HPDataFrame. If new.pix is specified then this function returns a new HPDataFrame with the same number of rows as x, but with pix attribute new.pix. Thus, new.pix must have length equal to nrow(x).

Usage

```
## S3 method for class 'HPDataFrame'
pix(x, new.pix, ...)
```

Arguments

x	a HPDataFrame .
new.pix	optional vector of pixel indices with length equal to nrow(x)
...	Unused arguments.

Value

The vector of HEALPix pixel indices (integers) or, if new.pix is specified, a new HPDataFrame.

Examples

```
df <- HPDataFrame(I = rep(0,12), nside = 1)
pix(df)
```

pix2coords	<i>Convert pixel indices to cartesian/spherical coordinates</i>
------------	---

Description

Convert HEALPix pixel indices to cartesian or spherical coordinates

Usage

```
pix2coords(nside, coords = "cartesian", ordering = "nested", spix)
```

pixelWindow	<i>Find high resolution pixels falling in a lower resolution window</i>
-------------	---

Description

Find all pixels in a higher resolution that fall within the specified pixel area at a lower resolution. All pixels are assumed to be in nested ordering.

Usage

```
pixelWindow(j1, j2, pix.j1)
```

Arguments

j1	An integer. The lower resolution, with $j1 \leq j2$. Note that <code>resolution = log2(nside)</code> .
j2	An integer. The upper resolution.
pix.j1	An integer. The pixel index at resolution j1 within which all pixels from resolution j2 will be returned. <code>pix.j1</code> can also be a vector of non-zero pixel indices.

Value

All pixels in resolution j2 that fall within the pixel `pix.j1` specified at resolution j1

Examples

```
pixelWindow(3, 3, 2)
pixelWindow(3, 4, 2)
pixelWindow(3, 5, 2)
```

plot.CMBCorrelation	<i>Plot sample CMBCorrelation</i>
---------------------	-----------------------------------

Description

Plots sample (empirical) correlation function. Uses [plot.variogram](#) from **geoR** package.

Usage

```
## S3 method for class 'CMBCorrelation'
plot(x, ...)
```

Arguments

x	An object of class CMBCorrelation.
...	Extra arguments as in plot.variogram passed to <code>plot.default</code> .

Value

Produces a plot with the sample correlation function.

References

geoR package, [corrCMB](#), [variog](#), [plot.variogram](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
corcmb <- corrCMB(cmbdf, max.dist = pi/2, num.bins = 6)
plot(corcmb)
```

plot.CMBCovariance	<i>Plot sample CMBCovariance</i>
--------------------	----------------------------------

Description

Plots sample (empirical) covariance function. Uses [plot.variogram](#) from **geoR** package.

Usage

```
## S3 method for class 'CMBCovariance'
plot(x, ...)
```

Arguments

x	An object of class CMBCovariance.
...	Extra arguments as in plot.variogram passed to plot.default.

Value

Produces a plot with the sample covariance function.

#' @references **geoR** package, [covCMB](#), [variog](#), [plot.variogram](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
Cov <- covCMB(cmbdf, max.dist = pi/2, num.bins = 6)
plot(Cov)
```

plot.CMBDataFrame	<i>Plot CMB Data</i>
-------------------	----------------------

Description

This function produces a plot from a [CMBDataFrame](#).

Usage

```
## S3 method for class 'CMBDataFrame'
plot(
  x,
  intensities = "I",
  add = FALSE,
  sample.size,
  type = "p",
  size = 1,
  box = FALSE,
  axes = FALSE,
  aspect = FALSE,
  col,
  back.col = "black",
  labels,
  hp.boundaries = 0,
  hpb.col = "gray",
  depth_test = "less",
  lab_depth_test = "always",
  ...
)
```

Arguments

x	A CMBDataFrame .
intensities	The name of a column that specifies CMB intensities. This is only used if col is unspecified.
add	If TRUE then this plot will be added to any existing plot. Note that if back.col (see below) is specified then a new plot window will be opened and add = TRUE will have no effect.
sample.size	Optionally specifies the size of a simple random sample to take before plotting. This can make the plot less computationally intensive.
type	A single character indicating the type of item to plot. Supported types are: 'p' for points, 's' for spheres, 'l' for lines, 'h' for line segments from z = 0, and 'n' for nothing.
size	The size of plotted points.
box	Whether to draw a box.

axes	Whether to draw axes.
aspect	Either a logical indicating whether to adjust the aspect ratio, or a new ratio.
col	Specify the colour(s) of the plotted points.
back.col	Optionally specifies the background colour of the plot. This argument is passed to <code>rgl::bg3d</code> .
labels	Optionally specify a vector of labels to plot, such as words or vertex indices. If this is specified then <code>rgl::text3d</code> is used instead of <code>rgl::plot3d</code> . Then <code>length(labels)</code> must equal <code>nrow(x)</code> .
hp.boundaries	Integer. If greater than 0 then HEALPix pixel boundaries at <code>nside = hp.boundaries</code> will be added to the plot.
hpb.col	Colour for the <code>hp.boundaries</code> .
depth_test	The depth test to be applied to the plotted points. This controls how resistant the plotted object is to being obscured. See rgl.material
lab_depth_test	The rgl depth test to be applied to the labels and pixel boundaries if present. See rgl.material
...	Arguments passed to <code>rgl::plot3d</code> .

Value

A plot of the CMB data

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
plot(cmbdf)
```

plot.CMBWindow	<i>Visualise a CMBWindow</i>
----------------	------------------------------

Description

Visualise a [CMBWindow](#)

Usage

```
## S3 method for class 'CMBWindow'
plot(
  x,
  add = TRUE,
  type = "l",
  col = "red",
  size = 2,
  box = FALSE,
```

```

    axes = FALSE,
    aspect = FALSE,
    back.col,
    depth_test = "always",
    ...
)

```

Arguments

<code>x</code>	A CMBWindow.
<code>add</code>	if TRUE then this plot will be added to any existing plot. Note that if <code>back.col</code> (see below) is specified then a new plot window will be opened and <code>add = TRUE</code> will have no effect
<code>type</code>	a single character indicating the type of item to plot. Supported types are: 'p' for points, 's' for spheres, 'l' for lines, 'h' for line segments from $z = 0$, and 'n' for nothing.
<code>col</code>	specify the colour(s) of the plotted points
<code>size</code>	the size of plotted points
<code>box</code>	whether to draw a box
<code>axes</code>	whether to draw axes
<code>aspect</code>	either a logical indicating whether to adjust the aspect ratio, or a new ratio.
<code>back.col</code>	specifies the background colour of the plot. This argument is passed to <code>rgl::bg3d</code> .
<code>depth_test</code>	The depth test to be applied. This controls how resistant the plotted object is to being obscured. See rgl.material
<code>...</code>	arguments passed to <code>rgl::plot3d</code>

Value

No meaningful return value; called for its side effect of drawing the window in the current `rgl` scene.

Examples

```

win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
win2 <- CMBWindow(theta = c(2*pi/3,3*pi/4,3*pi/4, 2*pi/3), phi = c(pi/4,pi/4,pi/3,pi/3))
plot(win1)
plot(win2)

```

plot.HPDataFrame	<i>Plot HPDataFrame</i>
------------------	-------------------------

Description

This function produces a plot from a [HPDataFrame](#). If columns x,y,z (cartesian) or theta,phi (colatitude and longitude respectively) are present in x, then these will be used as coordinates for plotting. Otherwise, the HEALPix indices as in pix(x) will be used. If HEALPix indices are used and multiple rows correspond to a single pixel index, then beware that values may be obfuscated in the plot, and all locations are pixel centers.

Usage

```
## S3 method for class 'HPDataFrame'
plot(
  x,
  intensities = "I",
  add = FALSE,
  sample.size,
  type = "p",
  size = 1,
  box = FALSE,
  axes = FALSE,
  aspect = FALSE,
  col = "blue",
  back.col = "black",
  labels,
  hp.boundaries = 0,
  hpb.col = "gray",
  depth_test = "less",
  lab_depth_test = "always",
  ...
)
```

Arguments

x	A HPDataFrame.
intensities	The column name for the data in x that is to be treated as intensities for plotting.
add	if TRUE then this plot will be added to any existing plot. Note that if back.col (see below) is specified then a new plot window will be opened and add = TRUE will have no effect
sample.size	optionally specifies the size of a simple random sample to take before plotting. This can make the plot less computationally intensive
type	a single character indicating the type of item to plot. Supported types are: 'p' for points, 's' for spheres, 'l' for lines, 'h' for line segments from z = 0, and 'n' for nothing.

size	the size of plotted points
box	whether to draw a box
axes	whether to draw axes
aspect	either a logical indicating whether to adjust the aspect ratio, or a new ratio.
col	specify the colour(s) of the plotted points
back.col	optionally specifies the background colour of the plot. This argument is passed to <code>rgl::bg3d</code> .
labels	optionally specify a vector of labels to plot, such as words or vertex indices. If this is specified then <code>rgl::text3d</code> is used instead of <code>rgl::plot3d</code> . Then <code>length(labels)</code> must equal <code>nrow(x)</code>
hp.boundaries	integer. If greater than 0 then HEALPix pixel boundaries at <code>nside = hp.boundaries</code> will be added to the plot
hpb.col	colour for the <code>hp.boundaries</code>
depth_test	The depth test to be applied to the plotted points. This controls how resistant the plotted object is to being obscured. This controls how resistant the plotted
lab_depth_test	The rgl depth test to be applied to the labels and pixel boundaries if present. See rgl.material
...	arguments passed to <code>rgl::plot3d</code>

Value

A plot of the data locations according to coordinate columns or HEALPix index

Examples

```
hpdf <- HPDataFrame(I = rep(0,12), nside = 1)
plot(hpdf, size = 5, col = "yellow", back.col = "black",
     hp.boundaries = 1)
```

plot.variogram	<i>Plot sample variogram</i>
----------------	------------------------------

Description

Plots sample (empirical) variogram. Uses [plot.variogram](#) from **geoR** package. See [plot.variogram](#) for usage and arguments.

Value

Produces a plot with the sample variogram.

References

geoR package, [variogramCMB](#), [variog](#), [plot.variogram](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
varcmb <- variogramCMB(cmbdf, max.dist = pi/2, num.bins = 6)
plot(varcmb)
```

plotAngDis	<i>Plot angular scatterplots and means</i>
------------	--

Description

For specified measurements from [CMBDataFrame](#) this function produces scatterplots and binned means versus theta and phi angles.

Usage

```
plotAngDis(cmbdf, intensities = "I")
```

Arguments

cmbdf	A CMBDataFrame object.
intensities	The name of a column of cmbdf, containing measured values.

Value

2x2 plot. The first row shows scatterplots. The second row gives barplots of the corresponding means computed over bins. The first column corresponds to the values of theta and the second one is for psi.

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
plotAngDis(cmbdf)
```

plotcovmodelCMB	<i>Plot theoretical CMB Covariance</i>
-----------------	--

Description

Plots theoretical covariance functions from the list defined in [covmodelCMB](#)

Usage

```
plotcovmodelCMB(
  cov.model = "matern",
  sigmasq = 1,
  phi = pi,
  kappa = 0.5,
  from = 0,
  to = pi,
  ...
)
```

Arguments

cov.model	A type of the correlation function. Available choices are: "matern", "exponential", "spherical", "powered.exponential", "cauchy", "gencauchy", "pure.nugget", "askey", "c2wendland", "c4wendland", "sinpower", "multiquadric". The default is "matern"
sigmasq	The variance parameter as documented in covmodelCMB . The default is 1.
phi	The range parameter as documented in covmodelCMB . The default is pi.
kappa	A smoothness parameter of the correlation function. The default is 0.5.
from	A lower range of the plotting region. The default is lb = 0
to	An upper range of the plotting region. The default is ub = pi.
...	optional plotting parameters.

Value

Produces a plot with the theoretical covariance function.

References

[covmodelCMB](#)

Examples

```
plotcovmodelCMB("matern", sigmasq = 5)
plotcovmodelCMB("askey", phi = pi/4, to = pi/2, kappa = 4)
```

plotvariogram	<i>Plot theoretical variogram</i>
---------------	-----------------------------------

Description

Plots theoretical variogram functions from the list defined in [covmodelCMB](#)

Usage

```
plotvariogram(
  cov.model = "matern",
  sigmasq = 1,
  phi = pi,
  kappa = 0.5,
  from = 0,
  to = pi,
  ...
)
```

Arguments

cov.model	A type of the variogram function. Available choices are: "matern", "exponential", "spherical", "powered.exponential", "cauchy", "gencauchy", "pure.nugget", "askey", "c2wendland", "c4wendland", "sinpower", "multiquadric". The default is "matern"
sigmasq	The variance parameter as documented in covmodelCMB . The default is 1.
phi	The range parameter as documented in covmodelCMB . The default is pi.
kappa	A smoothness parameter of the variogram function. The default is 0.5.
from	A lower range of the plotting region. The default is lb=0
to	An upper range of the plotting region. The default is ub= pi.
...	optional plotting parameters.

Value

Produces a plot with the theoretical variogram.

References

[covmodelCMB](#)

Examples

```
plotvariogram("matern", sigmasq = 5)
plotvariogram("askey", phi = pi/4, to = pi/2, kappa = 4)
```

practicalRangeCMB *Practical range for covariance function*

Description

This function computes the practical range for covariance functions on spheres. The function extends [practicalRange](#) from the package **geoR** to additional covariance models on spheres.

Usage

```
practicalRangeCMB(cov.model, phi, kappa = 0.5, correlation = 0.05, ...)
```

Arguments

cov.model	A type of the correlation function. Available choices are: "matern", "exponential", "spherical", "powered.exponential", "cauchy", "gencauchy", "pure.nugget", "askey", "c2wendland", "c4wendland", "sinpower", "multiquadric".
phi	The range parameter as documented in covmodelCMB
kappa	A smoothness parameter of the correlation function.
correlation	A correlation threshold (default is 0.05)
...	other optimisation parameters

Details

The practical(effective) range for a covariance function is the distance at which a covariance function first time reaches the specified value correlation. For covariance functions on spheres the practical range does not exceed π , the distance beyond which a covariance function is not defined. For the covariance functions "spherical", "askey", "c2wendland", "c4wendland" their practical ranges are equal to lengths of their support.

Value

Value of the practical range for the covariance function specified in [covmodelCMB](#)

References

geoR package, [practicalRange](#), [covmodelCMB](#)

Examples

```
practicalRangeCMB(cov.model = "sinpower", phi = 0.1, kappa = 0.5)
practicalRangeCMB(cov.model = "askey", phi = 0.1, kappa = 0.5)
```

print.CMBDataFrame	<i>Print CMBDataFrame</i>
--------------------	---------------------------

Description

This function neatly prints the contents of a CMBDataFrame.

Usage

```
## S3 method for class 'CMBDataFrame'  
print(x, ...)
```

Arguments

x	A CMBDataFrame .
...	arguments passed to print.tbl_df

Value

Prints contents of the CMB data frame to the console.

Examples

```
set.seed(1)  
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")  
print(cmbdf)
```

print.HPDataFrame	<i>Print a HPDataFrame</i>
-------------------	--

Description

This function neatly prints the contents of a HPDataFrame.

Usage

```
## S3 method for class 'HPDataFrame'  
print(x, ...)
```

Arguments

x	A HPDataFrame .
...	arguments passed to print.tbl_df

Value

Prints contents of the HPDataFrame to the console.

Examples

```
df <- HPDataFrame(I = rep(0,12), nside = 1, ordering = "nested")
print(df)
df
```

pwSpCorr

Power spectra estimate via correlation

Description

This function provides an angular power spectra estimate using the values of the sample correlations. The approach is based on Lawson-Hanson algorithm for non-negative least squares.

Usage

```
pwSpCorr(corcmb, lmax = 20 * length(corcmb$u))
```

Arguments

corcmb	An object of the class CMBCorrelation.
lmax	A number of angular power spectra components to estimate

Value

A data frame which first column is 1-d grid of l values from 0 to $l_{\max}$. The second column is estimated angular power spectra components on this grid.

References

Formula (2.1) in Baran A., Terdik G. Power spectrum estimation of spherical random fields based on covariances. *Annales Mathematicae et Informaticae* 44 (2015) pp. 15–22.

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
corcmb <- corrCMB(cmbdf, max.dist = pi, num.bins = 8)
pw <- pwSpCorr(corcmb, lmax = 10)
pw
```

qqnormWin	Normal QQ plot for CMBWindow
-----------	------------------------------

Description

This function is a modification of standard [qqnorm](#) functions to work with [CMBWindow](#) regions.

Usage

```
qqnormWin(cmbdf, win, intensities = "I")
```

Arguments

cmbdf	A CMBDataFrame .
win	A CMBWindow
intensities	A CMBDataFrame column with measured values.

Details

[qqnormWin](#) returns a normal QQ plot of for the specified [CMBDataFrame](#) column [intensities](#) and [CMBWindow](#) region. The function automatically adds a line of a “theoretical” normal quantile-quantile plot.

Value

A list with quantile components x and y and a normal QQ plot with QQ line

References

[qqnorm](#), [qqplot](#), [qqplotWin](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
win1 <- CMBWindow(theta = c(0, pi/2, pi/2), phi = c(0, 0, pi/2))
qqnormWin(cmbdf, win1)
```

qqplotWin

Quantile-Quantile plots for CMBWindows

Description

This function is a modification of standard [qqplot](#) functions to work with [CMBWindow](#) regions.

Usage

```
qqplotWin(cmbdf, win1, win2, intensities = "I")
```

Arguments

cmbdf	A CMBDataFrame .
win1	A CMBWindow
win2	A CMBWindow
intensities	A CMBDataFrame column with measured values.

Details

[qqplotWin](#) produces a QQ plot of quantiles of observations in two [CMBWindows](#) against each other for the specified [CMBDataFrame](#) column intensities. The function automatically adds a diagonal line.

Value

A list with quantile components x and y and a QQ plot with a diagonal line

References

[qqnormWin](#), [qqnorm](#), [qqplot](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
win1 <- CMBWindow(theta = c(0, pi/2, pi/2), phi = c(0, 0, pi/2))
win2 <- CMBWindow(theta = c(pi/2, pi, pi/2), phi = c(0, 0, pi/2))
qqplotWin(cmbdf, win1, win2)
```

qstat	<i>q-statistic</i>
-------	--------------------

Description

This function returns an estimated q-statistic for the specified column intensities in a [CMBDataFrame](#) and the list of [CMBWindows](#).

Usage

```
qstat(cmbdf, listwin, intensities = "I")
```

Arguments

cmbdf	A CMBDataFrame .
listwin	A list of CMBWindows
intensities	A CMBDataFrame column with measured values.

Details

The q-statistics is used to measure spatial stratified heterogeneity and takes values in [0, 1]. It gives the percent of the variance of intensities explained by the stratification. 0 corresponds to no spatial stratified heterogeneity, 1 to perfect spatial stratified heterogeneity.

Value

Estimated q-statistics for observations in a list of [CMBWindows](#)

References

Wang, J.F, Zhang, T.L, Fu, B.J. (2016). A measure of spatial stratified heterogeneity. Ecological Indicators. 67: 250–256.

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
win1 <- CMBWindow(theta = c(0, pi/2, pi/2), phi = c(0, 0, pi/2))
win2 <- CMBWindow(theta = c(pi/2, pi, pi/2), phi = c(0, 0, pi/2))
qstat(cmbdf, list(win1, win2))
```

rbind.CMBDataFrame [rbind](#) for CMBDataFrames

Description

Add a new row or rows to a [CMBDataFrame](#). All arguments passed to `...` must be [CMBDataFrames](#). If the [CMBDataFrame](#) arguments have overlapping pixel indices then all but one of the non-unique rows will be deleted unless `unsafe = TRUE`. If `unsafe = TRUE` then a [HPDataFrame](#) will be returned instead of a [CMBDataFrame](#).

Usage

```
## S3 method for class 'CMBDataFrame'
rbind(..., deparse.level = 1, unsafe = FALSE)
```

Arguments

<code>...</code>	A number of CMBDataFrames
<code>deparse.level</code>	See documentation for rbind.data.frame .
<code>unsafe</code>	A boolean. If the CMBDataFrame arguments have overlapping pixel indices then all but one of the non-unique rows will be deleted unless <code>unsafe = TRUE</code> . If <code>unsafe = TRUE</code> then a HPDataFrame will be returned instead of a CMBDataFrame .

Value

A [CMBDataFrame](#) containing the combined rows and windows, with overlapping pixel indices deduplicated. If `unsafe = TRUE`, returns an [HPDataFrame](#) retaining duplicate pixels instead. All inputs must have matching `nside` and `ordering`.

See Also

See the documentation for [rbind](#)

Examples

```
df <- CMBDataFrame(nside = 1, I = 1:12)

df.123 <- CMBDataFrame(df, spix = c(1,2,3))
df.123
df.234 <- CMBDataFrame(df, spix = c(2,3,4))
df.234

df.1234 <- rbind(df.123, df.234)
df.1234
class(df.1234) # A CMBDataFrame
pix(df.1234)

df.123234 <- rbind(df.123, df.234, unsafe = TRUE)
```

```
df.123234  
class(df.123234) # A HPDataFrame  
pix(df.123234)
```

rcosmo

Handling and Analysing CMB data

Description

Handling and Analysing Spherical, Healpix and Cosmic Microwave Background data on a HEALPix grid.

Details

The package `rcosmo` offers various tools for

- Downloading and transforming Cosmic Microwave Background radiation (CMB) and spherical data
- Working with Hierarchical Equal Area isoLatitude Pixelation of a sphere (Healpix)
- Spherical geometry
- Statistical analysis of CMB and spherical data
- Visualisation of Healpix data

Most of `rcosmo` features were developed for CMB, but it can also be used for other spherical data. It contains tools for transforming spherical data in cartesian and geographic coordinates to the Healpix representation.

Update

Current updates are available through URL: <https://github.com/frycast/rcosmo>

BugReports

<https://github.com/frycast/rcosmo/issues>

Author(s)

Daniel Fryer <d.fryer@latrobe.edu.au>, Andriy Olenko <a.olenko@latrobe.edu.au>, Ming Li <Ming.Li@latrobe.edu.au>, Yuguang Wang.

See Also

Useful links:

- <https://github.com/frycast/rcosmo>
- Report bugs at <https://github.com/frycast/rcosmo/issues>

resolution	<i>Get the arcmin resolution from a CMBDataFrame</i>
------------	--

Description

Get the arcmin resolution from a [CMBDataFrame](#)

Usage

```
resolution(cmbdf)
```

Arguments

cmbdf a [CMBDataFrame](#).

Value

The arcmin resolution as specified by the FITS file where the data was sourced

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
resolution(cmbdf)
```

ring2nest	<i>Convert ring to nest ordering.</i>
-----------	---------------------------------------

Description

ring2nest converts HEALPix pixel indices in the 'ring' ordering scheme to HEALPix pixel indices in the 'nested' ordering scheme.

Usage

```
ring2nest(nside, pix)
```

Arguments

nside is the HEALPix nside parameter (integer number 2^k)
 pix is a vector of HEALPix pixel indices, in the 'ring' ordering scheme.

Value

the output is a vector of HEALPix pixel indices in the 'nested' ordering scheme.

Examples

```
## Convert (1,2,23) from ring to nest at nside = 8
nside <- 8
pix <-c(1,2,23)
ring2nest(nside,pix)
```

sampleCMB

*Take a simple random sample from a [CMBDataFrame](#)***Description**

This function returns a [CMBDataFrame](#) which size equals to sample.size, whose rows comprise a simple random sample of the rows from the input CMBDataFrame.

Usage

```
sampleCMB(cmbdf, sample.size)
```

Arguments

cmbdf a [CMBDataFrame](#).
sample.size the desired sample size.

Value

A [CMBDataFrame](#) which size equals to sample.size, whose rows comprise a simple random sample of the rows from the input CMBDataFrame.

Examples

```
df <- CMBDataFrame(nside = 16, I = rnorm(12 * 16 ^ 2), ordering = "nested")
df.sample <- sampleCMB(df, sample.size = 100)
df.sample
```

siblings	<i>Return siblings of pixel</i>
----------	---------------------------------

Description

The siblings of pixel p are defined as the children of the parent of p . Note this is resolution independent.

Usage

```
siblings(p)
```

Arguments

p	Pixel index in nested order.
---	------------------------------

Value

An integer vector of four nested pixel indices sharing the same parent as p, including p itself.

Examples

```
siblings(11)
siblings(12)
```

sphericalHarmonics	<i>Compute spherical harmonic values at given points on the sphere.</i>
--------------------	---

Description

The function sphericalHarmonics computes the spherical harmonic values for the given 3D Cartesian coordinates.

Usage

```
sphericalHarmonics(L, m, xyz)
```

Arguments

L	The degree of spherical harmonic (L=0,1,2,...)
m	The order number of the degree-L spherical harmonic (m=-L,-L+1,...,L-1,L)
xyz	Dataframe for given points in 3D cartesian coordinates

Value

values of spherical harmonics

References

See https://en.wikipedia.org/wiki/Table_of_spherical_harmonics

It uses equation (7) in Hesse, K., Sloan, I. H., & Womersley, R. S. (2010). Numerical integration on the sphere. In Handbook of Geomathematics (pp. 1185-1219). Springer Berlin Heidelberg, but instead of the order $k=1,\dots,2L+1$ in the book we use $m=k-L-1$.

Examples

```
## Calculate spherical harmonic value at
## the point (0,1,0) with L=5, m=2
point<-data.frame(x=0,y=1,z=0)
sphericalHarmonics(5,2,point)

## Calculate spherical harmonic values at
## the point (1,0,0), (0,1,0), (0,0,1) with L=5, m=2
points<-data.frame(diag(3))
sphericalHarmonics(5,2,points)
```

summary.CMBDataFrame Summarise a [CMBDataFrame](#)

Description

This function produces a summary from a CMBDataFrame.

Usage

```
## S3 method for class 'CMBDataFrame'
summary(object, intensities = "I", ...)
```

Arguments

object	A CMBDataFrame .
intensities	the name of a column specifying CMB intensities (or potentially another numeric quantity of interest)
...	Unused arguments.

Value

A summary includes window's type and area, total area covered by observations, and main statistics for intensity values

Examples

```

ns <- 16
df <- CMBDataFrame(I = rnorm(12*ns^2), nside = ns,
                    ordering = "nested")

win1 <- CMBWindow(x=0,y=3/5,z=4/5,r=0.8)
df.sample1 <- window(df, new.window = win1)
summary(df)

```

summary.CMBWindow	Summarise a CMBWindow
-------------------	---------------------------------------

Description

This function produces a summary from a CMBWindow

Usage

```

## S3 method for class 'CMBWindow'
summary(object, ...)

```

Arguments

object	A CMBWindow.
...	Unused arguments.

Value

A summary includes window's type and area

Examples

```

win <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
summary(win)

win1<- CMBWindow(x=0,y=3/5,z=4/5,r=0.8, set.minus = TRUE)
summary(win1)

```

summary.HPDataFrame	Summarise a HPDataFrame
---------------------	---

Description

This function produces a summary from a HPDataFrame.

Usage

```
## S3 method for class 'HPDataFrame'
summary(object, intensities = "I", ...)
```

Arguments

object	A HPDataFrame.
intensities	the name of a column specifying intensities (or potentially another numeric quantity of interest)
...	Unused arguments.

Value

A summary includes window's type and area, total area covered by observations, and main statistics for intensity values

Examples

```
ns <- 2
hpdf <- HPDataFrame(I = rnorm(12*ns^2), nside = 2,
                     ordering = "nested")
win <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
hpdf.win <- window(hpdf, new.window = win)
summary(hpdf.win)
```

triangulate	Triangulate a polygonal CMBWindow
-------------	---

Description

Triangulate a polygonal [CMBWindow](#)

Usage

```
triangulate(win)
```

Arguments

`win` a CMBWindow object

Value

a list of CMBWindow polygons or `minus.polygons`, each having 3 vertices and representing a triangle. If `winType` of `win` does not include "minus" then these triangles have pairwise disjoint interiors and their union is equal to the original polygon, `win`. Otherwise, if `winType` of `win` does include "minus" the triangles are the same as for the non-minus type above, but have "minus" types.

Examples

```
## Example 1

win <- CMBWindow(theta = c(2*pi/3, 3*pi/4, 3*pi/4, 2*pi/3),
                  phi = c(pi/4, pi/4, pi/3, pi/3))

win
plot(win)
win1 <- triangulate(win)
win1
summary(win1[[1]])
plot(win1[[1]], add= FALSE, col="green")
plot(win1[[2]], col="blue")

## Example 2: triangulation minus-type polygon

win <- CMBWindow(theta = c(pi/5, pi/3, pi/4, pi/3, pi/5),
                  phi = c(pi/5, pi/5, pi/4, pi/3, pi/3), set.minus = TRUE)

win
plot(win)
summary(win)
win1 <- triangulate(win)
win1
plot(win1[[1]], add= FALSE, col="green")
plot(win1[[2]], col="blue")
plot(win1[[3]], col="yellow")
summary(win1[[1]])
summary(win1[[2]])
summary(win1[[3]])
```

variofitCMB

Estimates parameters of variograms

Description

This function estimates variogram parameters by fitting a parametric model from `covmodelCMB` to a sample variogram. The function extends `variofit` from the package **geoR** to additional covariance models on spheres.

Usage

```
variofitCMB(
  vario,
  ini.cov.pars,
  cov.model,
  fix.nugget = FALSE,
  nugget = 0,
  fix.kappa = TRUE,
  kappa = 0.5,
  simul.number = NULL,
  max.dist = vario$max.dist,
  weights,
  minimisation.function,
  limits = geoR::pars.limits(),
  messages,
  ...
)
```

Arguments

<code>vario</code>	An object of the class <code>variogram</code> obtained as an output of the function variogramCMB .
<code>ini.cov.pars</code>	A vector with initial values for the variogram parameters. The first parameter corresponds to the variance σ^2 . The second parameter corresponds to the range ϕ of the correlation function.
<code>cov.model</code>	A type of the variogram function. Available choices are: "matern", "exponential", "spherical", "powered.exponential", "cauchy", "gencauchy", "pure.nugget", "askey", "c2wendland", "c4wendland", "sinpower", "multiquadric". The default is "matern"
<code>fix.nugget</code>	logical. Indicates whether the nugget variance should be regarded as fixed or be estimated. The default is FALSE.
<code>nugget</code>	A value for the nugget parameter. Regarded as a fixed values if <code>fix.nugget = TRUE</code> or as a initial value for the minimization algorithm if <code>fix.nugget = FALSE</code> . The default is zero.
<code>fix.kappa</code>	logical. Indicates whether the parameter κ should be regarded as fixed or be estimated. The default is TRUE.
<code>kappa</code>	A value for the smoothness parameter. Regarded as a fixed values if <code>fix.kappa = TRUE</code> or as a initial value for the minimization algorithm if <code>fix.kappa = FALSE</code> . Required not in all covariance models, see covmodelCMB . The default is 0.5.
<code>simul.number</code>	number of simulation. Used if <code>vario</code> has empirical variograms for more than one data-set (simulations). The default is NULL
<code>max.dist</code>	A maximum distance to fit a variogram model. The default is <code>x\$max.dist</code> .
<code>weights</code>	Weights used in the loss function in the minimization algorithm.
<code>minimisation.function</code>	Minimization function ("optim", "nlm", "nls") to estimate the parameters.

limits	Lower and upper limits for the model parameters used in the numerical minimisation by <code>minimisation.function = "optim"</code> .
messages	logical. Indicates whether or not status messages are printed on the screen.
...	other minimisation parameters

Details

The parameter values of a variogram function from `covmodelCMB` are found by numerical optimization using one of the functions: `optim`, `nlm` and `nls`.

The function extends `variofit` from the package **geoR** to additional variogram models on spheres. Available models are: "matern", "exponential", "spherical", "powered.exponential", "cauchy", "gen-cauchy", "pure.nugget", "askey", "c2wendland", "c4wendland", "sinpower", "multiquadric".

Additionally it rescales an empirical variogram to the range $[0, 1]$ before numerical optimisation and then transforms all obtained results to the original scale. If `ini.cov.pars` are not provided then the 5×5 grid (`seq(0, max(vario$v), l=5)`, `seq(0, vario$max.dist, l=5)`) of initial values of σ^2 and ϕ is used.

Value

An object of the class `variomodel` and `variofit`, see `variofit`

References

geoR package, `variofit`, `covmodelCMB`

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
varcmb <- variogramCMB(cmbdf, max.dist = pi/2, num.bins = 6)
ols <- variofitCMB(varcmb, ini.cov.pars = c(1, 0.5),
                  cov.model = "matern", weights = "equal", messages = FALSE)
ols
```

variogramCMB

Sample variogram

Description

This function provides an empirical variogram for data in a `CMBDataFrame` or `data.frame`. It assumes that data are from a stationary spherical random field and the covariance depends only on a geodesic distance between locations. Output is a binned variogram.

Usage

```
variogramCMB(
  cmbdf,
  num.bins = 10,
  sample.size,
  max.dist = pi,
  breaks,
  equiareal = TRUE,
  calc.max.dist = FALSE
)
```

Arguments

<code>cmbdf</code>	is a CMBDataFrame or data.frame
<code>num.bins</code>	specifies the number of bins
<code>sample.size</code>	optionally specify the size of a simple random sample to take before calculating variogram. This may be useful if the full covariance computation is too slow.
<code>max.dist</code>	an optional number between 0 and pi specifying the maximum geodesic distance to use for calculating covariance. Only used if breaks are unspecified.
<code>breaks</code>	optionally specify the breaks manually using a vector giving the break points between cells. This vector has length <code>num.bins</code> since the last break point is taken as <code>max.dist</code> . If <code>equiareal = TRUE</code> then these breaks should be $\cos(r_i)$ where r_i are radii. If <code>equiareal = FALSE</code> then these breaks should be r_i .
<code>equiareal</code>	if TRUE then the bins have equal spherical area. If false then the bins have equal annular widths. Default is TRUE.
<code>calc.max.dist</code>	if TRUE then the <code>max.dist</code> will be calculated from the locations in <code>cmbdf</code> . Otherwise either <code>max.dist</code> must be specified or <code>max.dist</code> will default to pi.

Value

An object of class [variog](#) specified in the package **geoR**.

The attribute "breaks" contains the break points used to create bins. The result has `num.bins + 1` values since the first value at distance 0 is not counted as a bin.

u a vector with distances.

v a vector with estimated variogram values at distances given in **u**.

n number of pairs in each bin

sd standard deviation of the values in each bin

bins.lim limits defining the interval spanned by each bin

ind.bin a logical vector indicating whether the number of pairs in each bin is greater or equal to the value in the argument `pairs.min`

var.mark variance of the data

beta.ols parameters of the mean part of the model fitted by ordinary least squares

output.type echoes the option argument

max.dist maximum distance between pairs allowed in the variogram calculations
n.data number of data
direction direction for which the variogram was computed
call the function call

References

geoR package, [variog](#), [covCMB](#), [corrCMB](#)

Examples

```
set.seed(1)
cmbdf <- CMBDataFrame(nside = 8, I = rnorm(12 * 8^2), ordering = "nested")
varcmb <- variogramCMB(cmbdf, max.dist = pi/2, num.bins = 6)
varcmb
```

window

window generic

Description

Detailed descriptions and examples can be found in documentation for specific window functions [window.CMBDataFrame](#), [window.HPDataFrame](#), [window.data.frame](#)

Usage

```
window(x, ...)
```

Arguments

x	An object.
...	Extra arguments.

Value

The stored window (a CMBWindow, a list of windows, or NULL for full sky) when queried, or an object of the input class restricted to the requested sky region. See the individual methods.

See Also

[window.CMBDataFrame](#) [window.HPDataFrame](#) [window.data.frame](#)

window.CMBDat	<i>Get a sub window from a CMBDat object</i>
---------------	--

Description

This function returns a data.frame containing the data in `x` restricted to the CMBWindow `new.window`

Usage

```
## S3 method for class 'CMBDat'
window(x, new.window, intersect = TRUE, ...)
```

Arguments

<code>x</code>	a CMBDat object.
<code>new.window</code>	A single CMBWindow object or a list of them.
<code>intersect</code>	A boolean that determines the behaviour when <code>new.window</code> is a list containing BOTH regular type and "minus" type windows together (see details).
<code>...</code>	Unused arguments.

Details

Windows that are tagged with `set.minus` (see [CMBWindow](#)) are treated differently from other windows.

If the argument is a list of CMBWindows, then interiors of all windows whose `winType` does not include "minus" are united (let A be their union) and exteriors of all windows whose `winType` does include "minus" are intersected, (let B be their intersection). Then, provided that `intersect = TRUE` (the default), the returned data.frame will be the points of `cmbdat$data` in the intersection of A and B . Otherwise, if `intersect = FALSE`, the returned data.frame consists of the points of `x$data` in the union of A and B .

Note that if A (resp. B) is empty then the returned data.frame will be the points of x in B (resp. A).

Value

A CMBDataFrame containing the data in `x` restricted to the CMBWindow `new.window`

Examples

```
filename <- system.file("extdata", "CMB_testmap.fits", package = "rcosmo")
cmbdat <- CMBDat(filename)
win <- CMBWindow(x = 0, y = 0, z = 1, r = pi/2)
cmbdat.win <- window(cmbdat, new.window = win)
class(cmbdat.win)
```

window.CMBDataFrame *Get a sub window from CMBDataFrame*

Description

When new.window or in.pixels is unspecified this function returns the [CMBWindow](#) attribute of a CMBDataFrame. The return value is NULL if the window is full sky. When new.window is specified this function instead returns a new CMBDataFrame whose CMBWindow attribute is new.window

Usage

```
## S3 method for class 'CMBDataFrame'
window(x, new.window, intersect = TRUE, in.pixels, in.pixels.res = 0, ...)
```

Arguments

x	A CMBDataFrame .
new.window	Optionally specify a new window in which case a new CMBDataFrame is returned whose CMBWindow is new.window. new.window may also be a list (see details section and examples).
intersect	A boolean that determines the behaviour when new.window is a list containing BOTH regular type and "minus" type windows together (see details).
in.pixels	A vector of pixels at resolution in.pixels.res whose union contains the window(s) new.window entirely, or if new.window is unspecified then this whole pixel is returned.
in.pixels.res	An integer. Resolution (i.e., j such that $n_{\text{side}} = 2^j$) at which the in.pixels parameter is specified
...	Unused arguments.

Details

Windows that are tagged with set.minus (see [CMBWindow](#)) are treated differently from other windows.

If the argument new.window is a list of CMBWindows, then interiors of all windows whose winType does not include "minus" are united (let A be their union) and exteriors of all windows whose winType does include "minus" are intersected, (let B be their intersection). Then, provided that intersect = TRUE (the default), the returned CMBDataFrame will be the points of cmbdf in the intersection of A and B . Otherwise, if intersect = FALSE, the returned CMBDataFrame consists of the points of x in the union of A and B .

Note that if A (resp. B) is empty then the returned CMBDataFrame will be the points of x in B (resp. A).

Value

The window attribute of x or, if new.window/in.pixels is specified, a new CMBDataFrame.

Examples

```
## Example 1: Create a new CMBDataFrame with a window

cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian",
                      ordering = "nested")
win <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
cmbdf.win <- window(cmbdf, new.window = win)
plot(cmbdf.win)
window(cmbdf.win)

## Example 2: Change the window of an existing CMBDataFrame

cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian", ordering = "nested")
window(cmbdf) <- win2 <- CMBWindow(theta = c(pi/6,pi/3,pi/3, pi/6),
                                   phi = c(0,0,pi/6,pi/6))
plot(cmbdf)

## Example 3: union of windows

## Create 2 windows
win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
win2 <- CMBWindow(theta = c(2*pi/3,3*pi/4,3*pi/4, 2*pi/3),
                  phi = c(pi/4,pi/4,pi/3,pi/3))

plot(win1)
plot(win2)

## Create CMBDataFrame with points in the union of win1 and win2

cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian", ordering = "nested")
cmbdf.win <- window(cmbdf, new.window = list(win1, win2), intersect = FALSE)
plot(cmbdf.win)

## Example 4: intersection of windows

## Create 2 windows
win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
win2 <- CMBWindow(theta = c(pi/4,pi/3,pi/3, pi/4),
                  phi = c(pi/4,pi/4,pi/3,pi/3))

plot(win1)
plot(win2)

## Create CMBDataFrame with points in the intersection of win1 and win2

cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian", ordering = "nested")
cmbdf.win1 <- window(cmbdf, new.window = win1)
cmbdf.win12 <- window(cmbdf.win1, new.window = win2)
plot(cmbdf.win12)
plot(win1)
plot(win2)
```

```
## Example 5: intersection of windows with "minus" type

## Create 2 windows with "minus" type
win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2),
                  set.minus =TRUE)
win2 <- CMBWindow(theta = c(pi/4,pi/3,pi/3, pi/4),
                  phi = c(pi/4,pi/4,pi/3,pi/3),
                  set.minus =TRUE)

plot(win1)
plot(win2)

## Create CMBDataFrame with points in the intersection of win1 and win2

cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian", ordering = "nested")
cmbdf.win <- window(cmbdf, new.window = list(win1, win2))
plot(cmbdf.win)

## Example 6: intersection of windows with different types

##Create 2 windows, one with "minus" type

win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
win2 <- CMBWindow(theta = c(pi/4,pi/3,pi/3, pi/4),
                  phi = c(pi/4,pi/4,pi/3,pi/3),
                  set.minus =TRUE)

plot(win1)
plot(win2)

## Create CMBDataFrame with points in the intersection of win1 and win2

cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian", ordering = "nested")
cmbdf.win <- window(cmbdf, new.window = list(win1, win2), intersect = TRUE)
plot(cmbdf.win)

## Example 7: union of windows with different types

win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2), set.minus =TRUE)
win2 <- CMBWindow(theta = c(pi/4,pi/3,pi/3, pi/4), phi = c(pi/4,pi/4,pi/3,pi/3))
plot(win1)
plot(win2)

## Create CMBDataFrame with points in the union of win1 and win2

cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian", ordering = "nested")
cmbdf.win <- window(cmbdf, new.window = list(win1, win2), intersect = FALSE)
plot(cmbdf.win)
```

window.data.frame *Get a sub window from a data.frame*

Description

This function returns a data.frame containing the data in `x` restricted to the CMBWindow `new.window`

Usage

```
## S3 method for class 'data.frame'
window(x, new.window, intersect = TRUE, ...)
```

Arguments

<code>x</code>	A data.frame. Must have columns labelled x,y,z specifying cartesian coordinates, or columns labelled theta, phi specifying colatitude and longitude respectively.
<code>new.window</code>	A single CMBWindow object or a list of them.
<code>intersect</code>	A boolean that determines the behaviour when <code>new.window</code> is a list containing BOTH regular type and "minus" type windows together (see details).
<code>...</code>	Unused arguments.

Details

Windows that are tagged with `set.minus` (see [CMBWindow](#)) are treated differently from other windows.

If the argument is a list of CMBWindows, then interiors of all windows whose `winType` does not include "minus" are united (let A be their union) and exteriors of all windows whose `winType` does include "minus" are intersected, (let B be their intersection). Then, provided that `intersect = TRUE` (the default), the returned data.frame will be the points of x in the intersection of A and B . Otherwise, if `intersect = FALSE`, the returned data.frame consists of the points of x in the union of A and B .

Note that if A (resp. B) is empty then the returned data.frame will be the points of x in B (resp. A).

Value

A data.frame containing the data in `x` restricted to the CMBWindow `new.window`

Examples

```
win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))

cmbdf <- CMBDataFrame(inside = 4)
df2 <- coords(cmbdf, new.coords = "cartesian")
df <- as.data.frame(df2[,1:3])
df
df.win <- window(df, new.window = win1)
df.win
```

window.HPDataFrame *Get a sub window from a [HPDataFrame](#)*

Description

This function returns a HPDataFrame containing the data in hpdf restricted to the CMBWindow new.window. If the HPDataFrame has columns x,y,z or theta, phi then these will be used to determine locations with priority over the HEALPix indices in pix(hpdf) unless healpixCentered = TRUE is given. Note that if healpixCentered = TRUE then columns x,y,z or theta, phi will be discarded and replaced with pixel center locations.

Usage

```
## S3 method for class 'HPDataFrame'
window(x, new.window, intersect = TRUE, healpixCentered = FALSE, ...)
```

Arguments

x	A HPDataFrame .
new.window	Optional. A single CMBWindow object or a list of them.
intersect	A boolean that determines the behaviour when new.window is a list containing BOTH regular type and "minus" type windows together (see details).
healpixCentered	A boolean. If the HPDataFrame has columns x,y,z or theta, phi then these will be used to determine locations with priority over the HEALPix indices in pix(x) unless healpixCentered = TRUE is given. Note that if healpixCentered = TRUE then columns x,y,z or theta, phi will be discarded and replaced with pixel center locations.
...	Unused arguments.

Details

Windows that are tagged with set.minus (see [CMBWindow](#)) are treated differently from other windows.

If the argument is a list of CMBWindows, then interiors of all windows whose winType does not include "minus" are united (let A be their union) and exteriors of all windows whose winType does include "minus" are intersected, (let B be their intersection). Then, provided that intersect = TRUE (the default), the returned data.frame will be the points of df in the the intersection of A and B . Otherwise, if intersect = FALSE, the returned data.frame consists of the points of df in the union of A and B .

Note that if A (resp. B) is empty then the returned data.frame will be the points of df in B (resp. A).

Value

A HPDataFrame containing the data in x restricted to the CMBWindow new.window. Or, if new.window is unspecified, then the window attribute of x is returned instead (and may be NULL).

Examples

```

ns <- 16
hpdf <- HPDataFrame(nside = ns, I = 1:(12*ns^2))
hpdf

win1 <- CMBWindow(theta = c(0,pi/2,pi/2), phi = c(0,0,pi/2))
plot(hpdf); plot(win1)

hpdf.win <- window(hpdf, new.window = win1)
plot(hpdf.win, col = "yellow", size = 4, add = TRUE)
attributes(hpdf.win)
window(hpdf.win)
hpdf.win

```

winType*Get/change winType*

Description

Get/change the winType (polygon or disk) of a [CMBWindow](#). If new.type is missing then the winType of win is returned. Otherwise, a new window is returned with winType equal to new.type. If you want to change the winType of win directly, then use [winType<-](#), see the examples below.

Usage

```
winType(win, new.type)
```

Arguments

win	a CMBWindow object or a list of such
new.type	optionally specify a new type. Use this to change between "polygon" and "minus.polygon" or to change between "disc" and "minus.disc"

Value

If new.type is missing then the winType of win is returned. Otherwise a new window is returned with winType equal to new.type

Examples

```

win <- CMBWindow(theta = c(pi/2,pi/2,pi/3, pi/3), phi = c(0,pi/3,pi/3,0))
winType(win)

win1 <- CMBWindow(x=0,y=3/5,z=4/5,r=0.8)
winType(win1)
cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian",

```

```
                                ordering = "nested")
cmbdf.win1 <- window(cmbdf, new.window = win1)
plot(cmbdf.win1)

winType(win1) <- "minus.disc"
winType(win1)
cmbdf <- CMBDataFrame(nside = 64, coords = "cartesian",
                      ordering = "nested")
cmbdf.win1 <- window(cmbdf, new.window = win1)
plot(cmbdf.win1)
```

Index

ancestor, 4
areCompatibleCMBDFs, 5
as.CMBDataFrame, 6
assumedConvex, 7
assumedUniquePix, 8

baseNeighbours, 8

cbind, 9
cbind.CMBDataFrame, 9
chi2.empirical, 10
chi2CMB, 10
children, 11, 41, 51
CMBDat, 11, 91
CMBDataFrame, 5, 8–10, 12, 15, 19–21, 30–34, 36, 39, 40, 46, 48, 53, 55, 59, 61, 64, 69, 73, 75–78, 80, 81, 83, 88, 89, 92
CMBWindow, 7, 10, 13, 14, 16, 30–32, 37, 47, 65, 75–77, 84, 85, 91, 92, 95–97
coords, 15
coords.CMBDataFrame, 15, 15
coords.CMBWindow, 15, 16
coords.data.frame, 15, 17
coords.HPDataFrame, 15, 18
corrCMB, 19, 22, 63, 90
cov.spatial, 22–24
covCMB, 21, 21, 63, 90
covmodelCMB, 22, 46, 70–72, 86–88
covPwSp, 24
curve, 45

data.frame, 6, 12, 20, 21, 35, 39, 89
displayPixelBoundaries, 25
displayPixels, 26
download.file, 28
downloadCMBMap, 28
downloadCMBPS, 29

entropy, 30
entropyCMB, 30

exprob, 31
extrCMB, 32

fmf, 32
fRen, 33

geo2sph, 34
geoAngle, 35
geoArea, 36
geoArea.CMBDataFrame, 36, 36
geoArea.CMBWindow, 36, 37
geoArea.HPDataFrame, 36, 38
geoDist, 38

header, 39
healpixCentered, 40
HPDataFrame, 8, 12, 18, 19, 38, 40, 40, 45, 46, 48, 53, 56, 60, 61, 67, 73, 78, 85, 96

ibp2p, 42
is.CMBDat, 43
is.CMBDataFrame, 43
is.CMBWindow, 44
is.HPDataFrame, 45

lines.variomodel.variofit, 45, 46
linesCMB, 45

maxDist, 46
maxWindowDist, 47
minDist, 48
mmap, 11, 12

neighbours, 49
nest2ring, 50
nestSearch, 51
nlm, 88
nls, 88
nside, 52
nside.CMBDataFrame, 52, 53
nside.HPDataFrame, 52, 53

numeric2col, 54

optim, 88

ordering, 55

ordering.CMBDataFrame, 55, 55

ordering.HPDataFrame, 55, 56

p2bp, 57

p2ibp, 57

palette, 54

parent, 58

pix, 58

pix.CMBDataFrame, 58, 59, 59

pix.HPDataFrame, 58, 59, 60

pix2coords, 60

pixelArea, 61

pixelWindow, 62

plot.CMBCorrelation, 62

plot.CMBCovariance, 63

plot.CMBDataFrame, 64

plot.CMBWindow, 65

plot.HPDataFrame, 67

plot.variogram, 62, 63, 68, 68

plotAngDis, 69

plotcovmodelCMB, 70

plotvariogram, 71

practicalRange, 72

practicalRangeCMB, 72

print.CMBDataFrame, 73

print.HPDataFrame, 73

print.tbl_df, 73

pwSpCorr, 74

qqnorm, 75, 76

qqnormWin, 75, 75, 76

qqplot, 75, 76

qqplotWin, 75, 76, 76

qstat, 77

rbind, 78

rbind.CMBDataFrame, 78

rbind.data.frame, 78

rcosmo, 79, 79

rcosmo-package (rcosmo), 79

resolution, 80

rgl, 65, 68

rgl.material, 26, 65, 66, 68

ring2nest, 80

sampleCMB, 81

siblings, 82

sphericalHarmonics, 82

subWindow, 14

summary.CMBDataFrame, 83

summary.CMBWindow, 84

summary.HPDataFrame, 85

suppressMessages, 5

triangulate, 85

variofit, 86, 88

variofitCMB, 45, 46, 86

variog, 20–22, 63, 68, 89, 90

variogramCMB, 21, 22, 68, 87, 88

window, 90

window.CMBDat, 91

window.CMBDataFrame, 90, 92

window.data.frame, 90, 94

window.HPDataFrame, 90, 96

winType, 97